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(54) Title: MODEL FOR TESTING IMMUNOGENICITY OF PEPTIDES

(57) Abstract

Assay methods for determining whether a peptide is likely to be immunogenic are based on a computer modeling of binding to a Class II MHC DR1 receptor. This is confirmed by competitive inhibition binding assays. The peptides are useful for eliciting an immune response for vaccination or the production of antibodies or T-cells.

Applicants: Rina Aharoni et al.

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Exhibit 10

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MODEL FOR TESTING IMMUNOGENICITY OF PEPTIDES

Government Interest

The invention described herein may be manufactured, licensed and used by or for governmental purposes without the payment of any royalties to us thereon.

Cross Reference

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This application is a continuation-in-part of U.S. Patent application Serial No. 08/064,559, filed May 21, 1993, and the present application incorporates U.S. Patent Application Serial No. 08/064,559 in its entirety by reference.

15 Field of the Invention:

This invention relates to a means of predicting potential of a peptide for eliciting immune response.

Background of the Invention:

Among the numerous steps required for an immunological response to occur is the presentation of the antigen by macrophages to the B-cell or T-cell. This presentation is mediated by the Class I and Class II major histocompatibility complex (MHC) molecules on the surface of the cell. The MHC molecules hold antigens in the form of the peptide fragments and together with the receptor molecule on the T-cells, form a macromolecular complex that induces a response in the T-cell. Therefore, a necessary step in an immune response is the binding of the antigen to the MHC.

Recent single crystal X-ray structures of human and murine Class I MHC's have been reported. Analysis of these crystal structures have shown that antigenic peptides lie in the so-called binding cleft for presentation to the T-cell. This cleft is formed by α_1 and α_2 domains and by β -strands from each domain forming the floor. Furthermore, the sequence polymorphism among Class I molecules can result in alterations of the surface of the cleft forming different pockets. Peptide side chains may insert into these pockets. different pockets may interact with different side chains. This implies the mechanism for the peptide specificity of Class I MHC's. Peptides bound to the Class I MHC's in the crystal structures were found to have both the amino and There were few carboxy termini tightly held by the MHC. interactions near the middle of the cleft. Hence the bound peptide is allowed to bend slightly in the center. observed binding mode helped to explain the apparent partial specificity of peptide sequence and the allowed variation in peptide length found among peptides isolated from Class I MHC's.

The precise mode of binding of peptides to Class II MHC molecules is less clear. While a single crystal X-ray diffraction structure for the HLA-DR1 MHC has been shown, the coordinates have remained unavailable. However, currently available theoretical and experimental results help form a hypothesis that the binding of a peptide to Class II MHC is similar to that observed with Class I. First, it is noted that the Class II binding cleft is structurally similar to

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that of Class I. This was concluded based upon a sequence analysis of 26 Class I and 54 Class II amino acid sequences.

Unlike with Class I molecules, self-peptides isolated from murine I-Ab and I-Eb, from murine I-Ad and from human HLA-DR1 molecules were found to be varied in size (13 to 25 residues long). The peptides isolated from the murine I-Ab and I-Eb molecules had heterogenous carboxy termini while those from I-Ad and HLA-DR1 had ragged termini at both ends. The varying lengths indicate that the amino and carboxy termini of the peptides were not critical for the binding. One or both termini may protrude from the binding site and be available for further processing. The residues critical for binding were proposed to be at the ends of the peptide as opposed to the center.

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Summary of the Invention:

It is the purpose of this invention to provide a method for preliminary screening of peptides for ability to elicit an immune response. Structural homology techniques were used to model a receptor (the Class II MHC is exemplified). This model makes it possible to preliminarily screen peptides for antigenic properties. By modifying the peptide to "fit" into the receptor it is possible to identify methods of rendering non-immunogenic peptides immunogenic.

The preliminary screening of peptides for immunogenicity comprises the steps of (1) creating a molecular model of a receptor followed by minimizing the model created, 2) modeling a peptide to be tested and minimizing the model of the peptide, then testing the

fit of the model of the peptide into the model of the receptor to produce a composite minimized receptor/minimized peptide model.

Upon finding an acceptable fit, the peptide may then be screened by a binding assay for actual binding to Class II MHC as a further test for immunogenicity.

It has been found that when the model of the peptide can not be fitted into the model of the receptor, the peptide will lack immunogenicity. While not all peptide models which can be made to "fit" into to model of the receptor will be effective as immunogens, the screening methods of the invention may make it possible to avoid undue biological testing of inappropriate peptides. By using the model, it is also possible to alter peptides to accommodate the receptor. Hence, the invention has both predictive and drug design applications.

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Brief Description of the Figures:

Fig. 1 shows the HLA-aw68 α_1 and α_2 domains with DR1 α_1 and β_1 domains.

Figs. 2-30 are a printout of the minimized coordinates of the receptor.

Figs. 31 and 32 shows the effects of various peptides inhibiting the binding of labeled hemagglutinin in a competitive binding assay.

25 <u>Detailed Description of the Invention:</u>

In order to understand and better predict peptide interaction with Class II MHC's and as an aid for synthetic peptide vaccine design, a structural homology model of HLA-DR1 molecule was made

using the Class I HLA-aw68 as a reference molecule. For purposes of this analysis, numerous conserved residues were aligned leading to a proposed three-dimensional model for the Class II structure very similar to that of Class I. This model retained the overall conformation of a Class I MHC and agreed with a considerable amount of the published data. Furthermore, peptides shown to bind to DR1 were docked in the binding cleft of the model and analyzed. The results agree with the experimental binding data presented here. Hence, it is shown that the structural homology model reported here is useful for screening Class II MHC functionality.

It had been hypothesized that few peptide residues may be required for binding to DR1. By substituting residues into the influenza hemagglutinin 307-319 T-cell epitope (HA) it had been determined that a single tyrosine at 308 was required for binding. A synthetic peptide with the tyrosine at position 308 and a lysine at 315 was found to bind DR1 as well as the native peptide. Hence, it was concluded that few peptide residues determine the high affinity binding to DR1.

The peptides produced according to the present invention may be used alone or chemically bound to another peptide and/or carrier in order to elicit an immune response. An immune response is elicited by administering a peptide to an animal in an effective dose and by an effective route of administration. Typically the peptide will be administered with an immunologically acceptable carrier. The routes of administration, dosages, times between multiple administrations will be based on the particular peptide and are standard operations of those skilled in the art.

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Of particular interest are peptides from pathogenic microorganisms and neoplasms. In such an example, a vaccine may be formed with the peptide and any known immunological carrier and may be administered prophylactically or therapeutically. The immune response may be elicited for a number of reasons other than for prophylaxis or therapy such as increasing antibody production for the harvesting of antibodies, or increasing specific B-cell or T-cell concentration for the production of hybridomas or cellular therapy.

The choice of host animals is limited only to those capable of an immune response. Preferred hosts are mammals, more preferred are humans.

The vaccine may contain plural peptides with each peptide corresponding to the same or different antigens. The peptides may be used unbound or they may be chemically bound to another peptide or an unrelated protein or other molecule. A preferred vaccine preparation contains a plurality of peptides chemically bound to a larger more immunogenic peptide.

The peptide may be adsorbed, bound or encapsulated in a biodegradeable microsphere, microcapsule, larger carrier or a combination of these. The carrier may have a slow or controlled release property thereby releasing the peptide under appropriate conditions and times for enhanced immunization. This is particularly important when administering the peptide orally where stomach acid can degrade the peptide.

Another embodiment of the present invention is to modify the amino acid sequence of a peptide to enhance its immunogenicity.

This is done by modifying the natural peptide sequence to bind to

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the Class II MHC receptor DR1 with superior binding affinity for a Class II MHC receptor DR1 than the natural peptide sequence. This modified peptide is considered a synthetic peptide. Alternatively, the sequence may be modified to have a greater inhibition of HA (306-318) binding to a Class II MHC receptor DR1.

Many amino acid changes are acceptable in the formation of a synthetic peptide. The changes may be for similar types of amino acids such as leucine for isoleucine or they may be for diverse types such as tyrosine for lysine.

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Materials and Methods:

The structural homology model for the DR1 Class II MHC was constructed using the OUANTA molecular modeling package (vision 3.2, Molecular Simulations, Inc., Burlington, MA) with the CHARMM and Protein Design modules. After alignment of the sequences as described below, gaps and loops were energy minimized using 100 steps of steepest descents minimization followed by 100 steps of adopted basis set Newton-Rapheson (ABNR) minimization. were closed using a fragment database from a selected set of high-resolution crystal structures. The resulting structure was minimized in vacuo using 1000 steps of steepest descents followed by an additional 1000 steps of ABNR minimization. A distance related electrostatic function was used in all calculations with a dielectric constant of 1.0. Non-bound parameter lists were updated every 20 steps with a cutoff distance of 15.0Å. Non-bonded calculations were performed using a shifted potential function between 11.0Å and 14.0Å. An extended atom set was used with only

polar hydrogen atoms specifically placed. There were no explicit hydrogen bond energy calculations performed.

All peptides were initially modeled using QUANTA in an extended chain conformation and subjected to 500 steps of ABNR minimization. The resulting structures remained essentially in extended chain conformations. Individual peptides were manually docked in several different orientations into the binding cleft region of the minimized DR1 structure. The resulting bimolecular complex was subjected to 5000 steps of steepest descents minimization with non-bonded interactions updated every five steps. After minimization, bound peptides remained essentially in extended chain conformations. The lowest energy complexes for each peptide were selected for further analysis.

The selected peptide and DR1 complexes and the minimized DR1 model were subjected to the following molecular dynamics regimen: 300 steps of heating to 300°K, 600 steps of equilibration at 300°K, and 1100 steps of production dynamics. During this simulation, the DR1 C α atoms were constrained in their starting positions. All non-bonded interaction parameters were as stated for the minimization procedure. The lowest energy structure during the course of the production dynamics was selected and subjected to the 5000 step minimization procedure described previously with the C α restraints removed. The resulting structures were used for the binding energy calculations and for hydrogen bonding analysis.

Hydrogen bonds were determined using the QUANTA default parameters. Maximum allowed distances were 2.5Å between a hydrogen and the acceptor atom and 3.3Å between the donor and acceptor atoms.

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The minimum angle allowed between any set of atoms forming a hydrogen bond was 90°.

Competitive Inhibition Binding Assay:

Was labeled with ¹²⁵I. The labeled HA peptides were then allowed to interact with purified DR1 molecules during incubation to allow formation of peptide/DR1 complexes. After incubation, the peptide/DR1 composition was exposed to a native gel for chromatographic separation or passed through a spun column to separate labeled peptide/DR1 complex and free labelled peptide. When unlabeled peptides were added before incubation of labeled HA peptides and DR1, and if the unlabelled peptides had capacity for binding to DR1 simultaneous with ¹²⁵I-HA, there was a resultant decrease in radioactive signal associated with the DR1. The extent of this decrease directly related to the binding capacity of the unlabeled unknown peptide.

Structural Homology Model for the DR1 Molecule:

The structural homology model was created, the reference molecule being the crystal structure of HLA-aw68. The HLA-aw68 coordinates and subsequent sequence were obtained from the entry 2HLA in the Brookhaven Protein Data Bank released January 15, 1991, which is incorporated herein by reference. The sequence for the DR1 molecule was for the α_1 domain was reported by Klein and for the β_1 domain, the study reported by Todd et al. (Nature 329, 599 (1987)).

The sequence alignment is based on Brown et al. (Nature 332, 845 (1988)). The complete alignment and numbering scheme for both

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are seen in Figure 1. The Class II, $oldsymbol{eta}_1$ and Class I $lpha_2$ domains regions were conserved with some variations at the ends where the two MHC's have different loop regions. The fourth B-strand in the α_1 domain of HLA-aw68 (residues 30-38) is disrupted in the DR1 model. Only three residues are in a β -sheet conformation, probably due to the inserted glycine at position 28 before the strand and the large deletion in the loop region immediately after the strand. The two alpha-helical regions are clearly maintained. Both helices have been observed to be discontinuous in the Class I molecules and are similar in the DR1 model. The α_1 domain helix is long and curves from residues 49α to 76α without significant disruption. essentially a single continuous helix. However, the $lpha_2$ helical region is broken into two separate helices as with the Class I molecules. A short helix (52-63) is separated from a longer helix (68-94) by a deformed region without secondary structure. deformation is more pronounced in the DR1 model as opposed to the Class I molecules due to an insertion.

Influenza Hemagglutinin Peptide with DR1:

The amino acid residues 307-319 of influenza hemagglutinin (Pro-Lys- Tyr-Val-Lys-Gln-Asn-Thr-Leu-Lys-Leu-Ala-Thr) make up a well-documented linear T-cell epitope which has been shown to be HLA-DR1 restricted. With the demonstration that the influenza hemagglutinin epitope (referred to as the HA peptide) binds DR1, it was chosen to be modeled into the binding cleft.

The peptide was initially inserted into the cleft so that Leu 11 HA was in the vicinity of the hydrophobic pocket. This allowed Asn 7 to be near the middle charged and polar groups of the cleft.

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The remaining residue of the motif (Lys 2) was near the vicinity of the remaining charged and polar residues at the end of the cleft.

The only adjustment to the starting conformation was a slight rearrangement of the terminal peptide proline and Tyr 3 to alleviate obvious bad contacts.

After the energy minimization of the bimolecular complex, the total energy was reduced to 483 kcal/mol. This reduction in energy was accomplished by alleviation of several bad contacts and also be formation of several hydrogen bonds. The sticking feature of this mode is lack of hydrogen bonds in the carboxy terminal half of the peptide. Only one hydrogen bond is identified between the backbone carbonyl group of Leu 9 and the side chain of the β_1 Asn 77. In contrast, the amino terminal half has eleven identified interactions. Four of these interaction involve the peptide backbone residues Tyr 3, Val 4, and Gln 6. The remainder involve the side chains of Lys 2, Tyr 3, Lys 5 and Gln 6. Interestingly, Lys 5 is involved in more interactions (three) than Lys 2 (only 2). No interactions were observed as anticipated with Asn 7. Instead, it was the glutamine at position 6 donating a hydrogen bond to the α_1 Asn 62. No interactions were observed for the amino and carboxy termini.

HA-YK Peptide with DR1:

The binding of the HA-YK peptide (Ala-Ala-Tyr-Ala-Ala-Ala-Ala-Ala-Ala-Ala-Ala-Lys-Ala-Ala) to the DR1 model was tested. In aligning the peptide in the cleft, it was deemed logical to insert the tyrosine residue into the hydrophobic region of the binding cleft. The lysine would then be in position to interact with the

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hydrophilic groups in the other half of the cleft. The resulting peptide orientation is the opposite that used for the HA and the CS3 (defined below) peptides. With the peptide oriented as described, the final docking position for the peptide was unclear. The hydrophobic pocket is quite large, and, at least in this model, could accommodate the peptide tyrosine in a number of positions by sliding the peptide lengthwise through the cleft. However, repositioning the peptide also repositions the lysine. There were primarily two positions for the lysine: one with the lysine inside the cleft and the second with it outside. Of the two positions, the former was the lower in energy by 46 kcal/mol and had the greater number of interactions with the protein (11 vs. 7). Thus, the preferred orientation of the peptide appears to be with the lysine inside the binding cleft region.

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CS3 subunit Pilin Peptide with DR1:

The suspected T-cell epitope for CS3 pilus subunit 63-78

(Ser-Lys-Asn-Gly-Thr-Val-Thr-Trp-Ala-His-Glu-Thr-Asn-Asn-Ser-Ala)

was modeled with the DR1 molecule. The peptide was inserted with

lysine inside the cleft in the hydrophilic region. This placed the

Thr 5 in the center of the binding cleft and the tryptophane

(residue 8) near the hydrophobic region. The resulting minimized

model had ten interactions between the peptide and the protein,

three interactions with the peptide backbone and five with the

peptide side chains. The remaining two were with the amino terminal

of the peptide. All of the interactions were in either the first

three residues, His 10 or Glu 11 in the peptide. No interactions

were observed in the center of the cleft or residues four through nine.

CFA/1 with DR1:

A peptide identified as CFA/1 (colonization factor antigen)

(Val-Gly-Lys-Asn-Ile-Thr-Val-Thr-Ala-Ser-Val-Asp-Pro) was prepared

and an attempt was made to "fit" the molecule into the cleft of the

DR1. The lysine at position 3 prevented insertion of the peptide.

10 Results:

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The peptides chosen to dock in the DR1 model are shown in Table

1. The peptides were docked manually in several orientations into
the DR1 model. The peptides were then tested in biological binding
assays with the following results:

15 Table I

Peptide	Molecular Model predicted binding	Binding in the bioassay
HA (influenza hemagglutinin)	Yes	Yes
HA-YK (synthetic peptide)	Yes	Yes
CS3 Pilin subunit	Yes	Yes
CFA/1	No	No

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Quantitative measurement of the inhibition of CS3 63-78 and HA 306-318 as compared to controls is shown in Fig. 31.

The binding energy was calculated as the difference between the final DR1 and peptide complex and the sum of the energies for the minimized DR and peptide models individually. The data is shown in Table II.

Table II.

Peptide	Protein	Residues	Sequence	Binding Energy (kcal/mol)
НА	Influenza hemagglutinin	306-318	PKYVKQNTLKLAT	-283
на-үк	synthetic peptide		ААҮАААААКАА	-216
CS3	CS3 pilin subunit	63-78	SKNGTVTWAHETNNSA	-245

CS60 and CS6B with DR1

Colonization factor antigen IV (CFA/IV is an antigen on the surface of many enterotoxigenic *E. coli* one component of which is CS6. CS6 has two major subunits and a number of minor subunits. Several peptides from CS6 have been sequenced and assayed for potential inhibition of radiolabeled HA (306-318)/DR1 complex as a measure of immunogenicity. The sequences of the subunits are shown in Table III.

Table III.

Peptide	Amino Acid Residues	Sequence
CS6α6	63-75	DEYGLGRLVNTAD
CS6α7	80-92	IIYQIVDEKGKKK
CS6α8	111-123	LNYTSGEKKISPG
CS6ß1	3-15	WQYKSLDVNVNIE
CS6ß2	42-54	QLYTVEMTIPAGV
CS6ß3	112-124	TSYTFSAIYTGGE
CS6ß4	123-135	GEYPNSGYSSGTY
CS6ß5	133-145	GTYAGHLTVSFYS

These peptides were assayed for inhibition of radioactively labeled HA(306-318)/DR1. The results are demonstrated in Fig. 32.

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The foregoing description of the specific embodiments reveal the general nature of the invention so that others can, by applying current knowledge, readily modify and/or adapt for various applications such specific embodiments without departing from the generic concept, and, therefore, such adaptations and modifications should and are intended to be comprehended within the meaning and range of equivalents of the disclosed embodiments. It is to be understood that the phraseology or terminology employed herein is for the purpose of description and not of limitation.

All references mentioned in this application are incorporated by reference.

We Claim:

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1. A method of preliminarily screening peptides for immunogenicity comprising the steps of:

- 1) creating a molecular model of receptor DR1 Class II MHC and minimizing the model of the DR1;
 - 2) modeling a peptide to be tested and minimizing the model of the peptide; and
- 3) testing fit of model obtained in step 2 into the model
 10 obtained in step 1 to produce a composite receptor/peptide model.
 - 2. A computerized model comprising a model of the DR1 molecule having fitted in a cleft therein a model of a peptide.
- 3. A method of claim 1 wherein, additionally, the receptor/peptide model is subjected to computer-simulated heating.
 - 4. A method of claim 1 further comprising, assaying the peptide by competitive inhibition binding to a Class II MHC receptor DR1.
 - 5. A minimized peptide capable of binding to a Class II MHC receptor DR1 and inhibiting the binding of HA (306-318).
- 6. A synthetic peptide, wherein the amino acid sequence of the
 minimized peptide according to claim 5 has been modified to have a
 superior binding affinity for a Class II MHC receptor DR1 to form at
 least a portion of the synthetic peptide.

7. A synthetic peptide, wherein the amino acid sequence of the minimized peptide according to claim 5, has been modified to have greater inhibition of HA (306-318) binding to a Class II MHC receptor DR1 to form at least a portion of the synthetic peptide.

- 8. A synthetic peptide according to claim 6, wherein an amino acid has been modified from a charged amino acid to an uncharged amino acid.
- 9. A synthetic peptide according to claim 7, wherein an amino acid has been modified from a charged amino acid to an uncharged amino acid.
- 10. A synthetic peptide according to claim 8, wherein saiduncharged amino acid is alanine.
 - 11. A synthetic peptide according to claim 9, wherein said uncharged amino acid is alanine.
- 12. A minimized peptide according to claim 5, wherein the sequence is selected from the group consisting of PKYVKQNTLKLAT, AAYAAAAAAKAA and SKNGTVTWAHETNNSA.
- 13. A minimized peptide according to claim 5, wherein the sequence is contained in a CFA.

14. A minimized peptide according to claim 13, wherein the sequence is selected from the group consisting of DEYGLGRLVNTAD, IIYQIVDEKGKKK, LNYTSGEKKISPG, WQYKSLDVNVNIE, QLYTVEMTIPAGV, TSYTFSAIYTGGE, GEYPNSGYSSGTY and GTYAGHLTVSFYS.

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- 15. A vaccine comprising:
 - a minimized peptide according to claim 5; and an immunologically acceptable carrier.
- 10 16. A vaccine comprising:
 - a synthetic peptide according to claim 6; and an immunologically acceptable carrier.
 - 17. A vaccine comprising:
- a synthetic peptide according to claim 7; and an immunologically acceptable carrier.
 - 18. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 15.
 - 19. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 16.

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20. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 17.

av68 DR1 #,	1 GSUSHRYF 9 YISVSUPGRG 19 EPRPIAVGYV 29 D DIGFVRPD 38 SOAASQRHEP 48 RAPHIEQEGP 1 GSUSHRYF 9 YISVSUPGRG 19 EPRPIAVGYV 29 D DIGFVRPD 38 SOAASQRHEP 48 RAPHIEQEGP 1 KEEUVIIQA 11 EFYLN PDQ 19 SG EFHFDP 27 DGDEIFHVDH 37 AXX 40 ETVARLEEPPG 1 IKEEUVIIQA 11 EFYLN PDQ 19 SG EFHFDP 27 DGDEIFHVDH 37
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3008	172 LENGKLILUK
DR1 .	
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| Conserved residues

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249	25 ASP 0	-48.57317 -29.35714 102.56241 A1	25	0.00000
250	26 PHE N	-50.15819 -29.70261 101.00377 A1	2 6	0.00000
251	26 PHE H	-50.60050 -29.38662 100.15349 A1	26	0.00000
252	26 PHZ CA	-50.56030 -30.99701 101.54898 Al	26	0.00000

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253	26 PHE	Св	51.4237	8 -30.8309	8 102.82009) A1	26	0.00000
254	26 PHZ	CG		6 -29.9203			26	0.00000
255	26 PHE			B -28.5567		λl	26	0.00000
256	26 PHE			0 -30.42265			26	0.00000
257	26 PHE			7 -27.6870			26	0.00000
258	26 PHE			4 -29.55076 9 -28.18547			26 26	0.00000
259	26 PHE 26 PHE			0 -31.82500			26	0.00000
260 261	26 PHE			7 -31.38211			26	0.00000
262	27 ASP			0 -33.08225			27	0.00000
263	27 ASP			8 -33.34660			27	0.00000
264	27 ASP			5 -34.11623			27	0.00000
265	27 ASP	CB		1 -34.25359			27	0.00000
266	27 ASP			5 -35.66101			27	0.00000
267	27 ASP	ODl	-52.95640	0 -36.23164	101.40272		27	0.00000
268	27 ASP	OD2		2 -36.19966			27 27	0.00000
269 270	27 ASP 27 ASP	0		2 - 34.05698 5 - 34.10386			27	0.00000
271	28 GLY	и		5 -33.92246			28	0.00000
272	28 GLY		49.78144				28	0.00000
273	28 GLY	CA		-33.86633			28	0.00000
274	28 GLY	c		-32.51821			28	0.00000
275	28 GLX	O		.,-32.03342		Al	28	0.00000
276	29 ASP	N		31.96145		A1	29	0.00000
277	29 ASP	н	-51.93955	-32.39453	96.87595	λl	29	0.00000
278	29 ASP	СА	-51.52130	-30.85124	95.43198		29	0.00000
279	29 ASP	CB.	-52.38232		94.30287	Al	29	0.00000
280	29 ASP	CG		-30.54960			29	0.00000
281	29 ASP	OD1		-30.32445	92.44162		29 29	0.00000
282 283	29 ASP 29 ASP	OD2		-30.10545 -29.64631	92.75048		29	0.00000
284	29 ASP	C O		-28.59407	95.46732		29	0.00000
285	30 GLU	n		-29.81701	97.32405		30	0.00000
286	30 GLÜ	н		-30.57647	97.90251		30	0.00000
287	30 GLU	CA		-28.70871	97.86160	Al	30	0.00000
288	30 GLÜ	CB	-54.71008	-29.18152	98.54724	Al	30	0.00000
289	30 GLU	CG	-55.84100		97.67444	Al	30	0.00000
290	30 CFA	CD		-29.50058	98.43828		30	0.00000
291	30 GLU	OE1		-30.47326	98.78969		30 30	0.00000 0.00000
292	30 GLU	OE2	· ·	-28.34297 -27.80995	98.66447 98.82987		30	0.00000
293 294	30 GFA 30 GFA	С 0	-52.13191	-28.20198	99.84476		30	0.00000
295	31 ILE	и	-52.76110	-26.53268	98.46810		31	0.00000
296	31 ILE	. н	-53.28532	-26.29454	97.65353		31	0.00000
297	31 ILE	Cλ	-51.98442	-25.52556	99.18511	አ1	31	0.0000
298	31 ILE	CB	-51.81933	-24.32331	98.23317	A.1	31	0.0000
299	31 ILE	CG2		-23.65625	97.91607		31	0.00000
300	31 ILE	CG1	-50,75168		98.71478		31	0.00000
301	31 ILE	CD	-50.41981		97.66271		31	0.00000
302	31 ILE	C		-25.12446 -24.97567			31 31	0.00000 0.00000
303 304	31 ILE 32 PHE	N.		-24.95576			32	0.00000
305	32 PHE	н.	-50 59061	-25.06998	101.18020	λ1	32	0.00000
306	32 PHE	CA	-51.85606	-24.57978	102.84494	Al	32	0.00000
307	32 PHE	CB	-50.76201	-25.18703	103.73391	Al	32	0.00000
308	32 PHE	CG	-51.19588	-26.22218	104.74893	λl	32	0.00000
309	32 PHE	CD1	-50.23927	-27.18148	105.15120	A.l	32	0.00000
310	32 PHE	CD2	-52.49149	-26.22907	105.32111		32	0.00000
311	32 PHS	CEl	-50.57386	-20.14516	106.12524	Al 11	32 32	0.00000 0.00000
312	32 PHE	CE2	-52.82688	-219600	106 60566	V]	32	0.00000
313 314	32 PHE	CZ C	-51.86506	-28.15152 -23.07181	103.07985		32	0.00000
314	32 PHE 32 PHE	0	-52.76581	-23.07101	103.62205	λl	32	0.00000
315	32 PRE	N	-50.69098	-22.53451	102.70813	ΑĴ	33	0.00000
	JJ 8.23	••	20.02953					

FIG. 6

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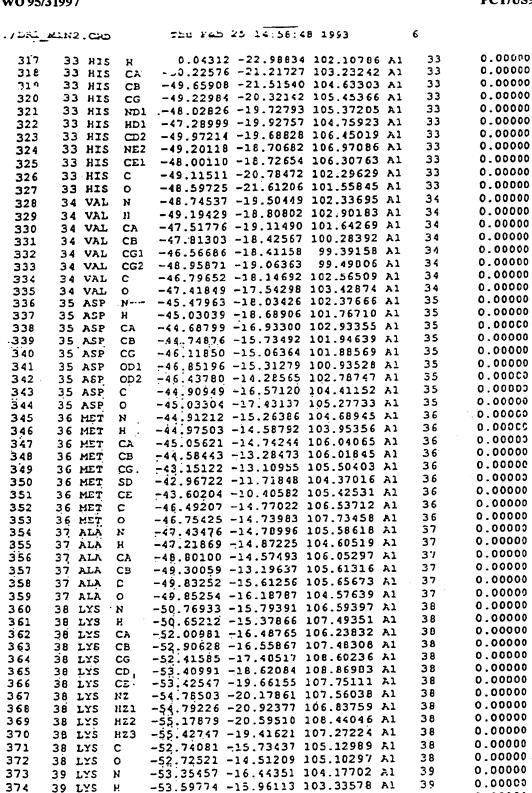


FIG. 7

-53.67982 -17.86931 104.22041 Al

-55.18971 -18.00506 104.55346 Al

-56.24681 -17.84030 103.42417 A1

-56.27039 -15.52580 102.62130 A1

-56.64532 -16.69161 101.13213 Al

-55,66137 -17.54082 100.43802 A1

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381	39 LYS	нгі	55.9564	5 -18.5458	1 100.36313	A1	3	0.00000
382	39 LYS						39	0.00000
3 6 3	39 LYS	нгз		6 -17.5192		A1	39	0.00000
384	39 LYS	С		3 -18.4428		λl	39	0.00000
385	39 LYS	0	-53.4335	1 -17.6751			39	0.00000
386	40 GLU	N	-53.4133				40	0.00000
387	40 GLU			4 -20.4369			40	0.00000
388	40 GLU			2 -20.14649			40	0.00000
389	40 GLU			2 -20.84643			40	0.00000
390	40 GLU			3 -19.82757			40 40	0.00000
391	40 GLU			7 -18.90107 1 -17.68372			40	0.00000
392	40 GLU 40 GLU			6 -19.35560			40	0.00000
393 394	40 GLU	C C		5 -20.74566			40	0.00000
395	40 GLU	0		1 -20.18304			40	0.00000
396	41 THR	N		1 -21.84976			41	0.00000
397	41 THR	11		-22.20472			41	0.00000
398	41 THR	CA		5 -22.55464		٨l	41	0.00000
399	41 THR	CB		-24.05830		Al	41	0.00000
400	41. THR	· OG1		2 -24.75407			41	0.00000
401	41 THR	HG1	- •	-25.59614			41	0.00000
402	41 THR	CG2	-55.95304				41	0.00000
.403	41 THR	C	-57.85722				41 41	0.00000
404	41 THR	0	-57.54877				42	0.00000
405	42 VAL	N H	-59.12316	5 -22.41525 -22.81422			42	0.00000
406 407	42 VAL 42 VAL	CA	-60.29134				42	0.00000
408	42 VAL	CB	-61.57611				42	0.00000
409	42 YAL	CG1		-21.83994	•		42	0.00000
410	42 VAL	CG2	-61.49852				42	0.00000
411	42 VAL	Ç			105.14170	Al.	42	0.00000
412	42 YAL	O	-60.54018	-22.16641	106.20404	Al	42	0.00000
413	43 TRP	N ·	-60.26652	-24.07997			43	0.00000
414	43 TRP	н.		-24.53437			43	0.00000
415	43 TRP	CA	-60.35178				43	0.00000
416	43 TRP	CB.	-60.91225				43	0.00000
417	43 TRP	CG	-60.96314				43 43	0.00000
418	43 TRP 43 TRP	CD2	-59.88712	-27.12198 -27.49629			43	0.00000
419 420	43 TRP 43 TRP	CE2 CE3	-60.47947 -58.50029	-27.24274	104.12574		43	0.00000
421	43 TRP	CD1		-26.86040	103.93459		43	0.00000
422	43 TRP	NEI	-61.84174	-27.31306	102.67766		43	0.00000
423	43 TRP	HE1	-62.50030	-27.49046	101.97106	Al	43	0.00000
424	43 TRP	CZ2	-59.62565	-27.96652	101.58525		43	0.00000
425	43 TRP	CZ3	-57.69263	-27.72260	103.07694	λl	43	0.00000
426	43 TRP	CH2	-58.24204	-28.07937	101.82597	Y1	43 43	0.00000 0.00000
427	43 TRP	C	-59.07150	-24.84373	107.16564	F 3	43	0.00000
428 429	43 TRP	0	-58.85050	-25.080/3	108.02836	ומ רא	44	0.00000
430	44 ARG	N H	-50.22471	-23.04303	106.07364	Al	44	0.00000
431	44 ARG	CA	-57 22046	-23.27577	107.89144	7.1 7.1	44	0.00000
432	44 ARG	CB		-23.69727		Al	44	0.00000
433	44 ARG	CG			108.22026		44	0.00000
434	44 ARG	CD	-53.30915	-23.64273	107.47006	A1	44	0.00000
435	44 ARG	NE	-52.18625	-23.26000	108.32672	Al	44 .	0.00000
436	44 ARG	HE	-52.36221	-22.91543	109.25139	λl	44	0.00000
437	44 ARG	CZ	-50.93149	-23.33650	107.86475		4.4	0.00000
438	44 ARG	נאא	-49.92358	-22.86519	108.56844	Al	4.4	0.00000 0.00000
439	44 ARG	LINH	-48.98076	-22.84912	108.24673		4 4 4 4	0.00000
440	44 ARG		-50.05162	-22.32230	109.51721	A.1	44	0.00000
441 442		КН2 НН21		-23.55024	106.89332	 A.1	44	0.00000
443			-51.41769		106.14195		14	0.00000
444		C		-22.05857			44	0.00000

FIG. 8



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445	44 ARC	. 0	6.5306	4 -21.3920	1 108.88943	A1 44	0.00000
446	45 LEU		-58.6527	0 -21.5568	6 108.18548	A1 45	
447	45 LEU		-59.3684	6 -22.0638	2 107.69729	A1 45	0.00000
448	45 LEU		-58.9467	9 -20.2189	7 108.70611 5 107.74942	λ1 45 λ1 45	0.00000
449 450	45 LEU 45 LEU		:-59,8726	7 -19.4631.	0 106.49579	A1 45	0.00000
451	45 LEU		-59.1303	4 -18.4190	6 105.47407	λ1 45	0.00000
452	45 LEU		-58:1043	1 -17.9228	9 106.87944	A1 45	0.0000
453	45. LEU		-59.5581	8 -20.2010	4 110.09168	λ1 45	0.00000
454	45 LEU	0	-59.5336	2 -19.1992	6 110.79199	A1 45	0.00000
455	46 GLU		-60.0891	7 -21.3602	4 110.48954	A1 46	0.00000
456	46 GLU		-60,1404	7 -22.14700	8 109.87631 7 111.86481	A1 46	0.00000
457 458	46 GLU 46 GLU		-61.4781	7 -22.7151 <i>1</i>	3 111.95437	X1 46	0.00000
459	46 GLU	CG	-60.8088	1 -24.0238	5 111.51772	A1 46	0.00000
460	46 GLU	CD	-61.8526	4 -25.11830	111.46594	A1 46	0.00000
461	46 GLU	OE1	-62.2686	7 -25.47222	110.36424	A1 46	0.00000
462	46 GLU	OE2			112.52280		0.00000
463 464	46 GLU 46 GLU	.o			114,10493		0.00000
465	47 GLU	H	-58.27888	-21.79310	112.44220	A1 47	0.00000
466	47 GLU	н	-58.18142	-22.06659	111.48920	Al 47	0.00000
.467	47 GLU	CA	-570815	-21.73864	113.27698	A1 47	0.00000
468	47 GLU	CB	-55.89123	-22.09481	112.39059	ሕ1 47	0.00000
469	47 GLU	CG	-55.95036	-23.50661	111.78977	ሕ1 47	0.00000
470	47 GLU	CD	-55.55418	-24.58648	112.78701	A1 47	0.00000
471 472	47 GLU 47 GLU	OE2			112.37878		0.00000
473	47 GLU	C	-56 A3A27	-24.27070 -20.38151	113.91510	A1 47	0.00000
574	47 GLU	o	-56.67000	-19.35742	113.26194		0.00000
475	48 PHE	N	-56.83807	-20.40716	115.24713	ሕ 1 48	0.00000
476	48 PHE	P.	-56.92233	-21.28123	115.72305	አ1 48	0.00000
477	48 PHE	CA	-56.75617	-19.13647	115.96362	A1 48 A1 48	0.00000
478	48 PXE	CB	-57.20231	_19.34982 _19.49841	117.41420 117.70829		0.00000
479 480	48 PHE	CG CD1	-59.71122	-19.00320	117.44215	A1 48	0.00000
481	48 PHE	CD2	-58.25459	-17.19966	118.24455	<u>A1 48</u>	0.00000
482	48 PHE	CE1	-60.84630	-18.20670	117.71246	A1 48	0.00000
483	48 PHE	CE2		-16.40286	118.51469	A1 48	0.00000
484	48 PHE	CZ	-60.68202	-16.908/8	118.24779	Al 48 Al 48	0.00000
485 486	48 PHE 48 PHE	C O	-55.39240 -54.35117	-10.46628	115.93777 115.72587	A1 48	0.00000
487	49 GLY	N	-55.43214	-17.15361	116.19781	A1 49	0.00000
488	49 GLY	В	-56.32069	-16.72101	116.34575	Al 49	0.00000
489	49 GLY	CV	-54.23420	-16.30520	116.13162	Al 49	0.00000
490	49 GLY	С	-52.92695	-16.86464	116.68056	A1 49 A1 49	0.00000 0.00000
491 492	49 GLY 50 ARG	0	-51.85815	-10.69886	116.10714	A1 50	0.00000
492	50 ARG	N H	-53.03313	-17.67284	118.24069	A1 50	0.00000
494	50 ARG	CY	-51.84656	-16.16800	118.42224	A1 50	0.00000
495	50 ARG	CB.	-52.27437	-18.92760	119.68340	al 50	0.00000
496	50 ARG	CG	-51.10260	-19.32736	120.58254	A1 50	0.00000
497	50 ARG	CD	-53.53002	-20.14774	121.80047	Al 50 Al 50	0.00000
498	50 ARG	ИE	-50.37432	-2U.44943	122.64580	A1 50	0.00000
499 500	50 ARG 50 ARG	HE CZ	-50.24449	-21.63184	123.26571	A1 50	0.00000
501	50 ARG	NHI	-49.18578	-21.84466	124.04503	እነ 50	0.0000
502	50 ARG	HH11	-49.05579	-22.71179	124.52612	A1 50	0.00000
503	SO ARG	нн12	-48.49360	-21.13233	124.16453	A1 50	0.00000
504	50 ARG	NE2	-51.15926	-22.58629	123.10823	h) 50 h) 50	0.00000
505	50 ARG	HH21	-51.08073	-23.47414	123.56513 A	A) 50 A) 50	0.00000
506 507	50 ARG 50 ARG	HH22 C	-51.06703	-19.09561	117.49074	A1 50	0.00000
508	50 ARG	0	-49.84240	-19.09133	117.41926	A: 50	0.00000
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509	51 PHE	ห	. 1.8491	5 -19.8777	8 116.73839) A1	5 i	0.00000
510	51 PHE				7 116.76797		51	0.00000
513	SI PHE	• • •	-		6 115.73402		51	0.00000
512	51 PHE				3 115.15449		51	0.00000
513	51 PHE				4 114.02359		51	0.00000
514	51 PHE		-51 8788I	-23.8285	1 114.29503	B Al	51	0.00000
515	51 PHE		-52 45781	-22.06484	112.69289	λ1	51	0.00000
516	51 PHE	CEI	-51 72481	-24.7442	1 113.22965) A1	51	0.00000
517	51 PHE		-52 30778	-22.98082	2 111.63059	A1	51	0.00000
	51 PHE	CZ	-51 04314	-24.31860	111.90000	λ1	51	0.00000
518	51 PHE	C	-51.54514 -EA 50269	-20 00696	114.69371	λ1	51	0.00000
519	51 PHE		-40.30260	-20.0000	114.47961	Al	51	0.00000
520		0	-45.31201	-10.2000	114.10202	A1	52	0.00000
521	52 ALA	.N	-51.20403	-19.05212	114.33870	1.3	52	0.00000
522	52 ALA	H	-50 54804	-10.05/04 -18 16484	113.11957	Al	52	0.00000
523	52 ALA	CA	-50.54696	-10.10400	112.62511	A I	52	0.00000
524	52 ALA	CB	-31.32030	-17.03210	113.62687	A 1	52	0.00000
525	52 ALA	C	49.28237	17 20034	112.94510	וג	52	0.00000
526	52 ALA	0	-40.27000	-17.03059	114.88764	A 1	53	0.00000
527	53 SER	,N	-49.35763 	-17.04933	114.00/03	7.1	53	0.00000
528	53 SER			-1/.11770	115.53815	N.	53	0.00000
529	53 SER	CA					53	0.00000
530	53 SER	CB	-48.58146	-16.02158	116.95383	VI	53	0.00000
531	53 SER	OG	-4759593	-15.15541	117.52845	WI.	53	0.00000
532	53 SER	HG	-47.85158	-14.90357	118.42105) 1	53	0.00000
533	53 SER	C	-46.99433	-17.42737	115.57426	7.1 7.1	53	0.00000
531	53 SER	0	-45.89463	-17.11790	115.12568	Y 7	54	0.00000
535	54 PHE	н.	-47.26082	-18.64220	116.08200	V.1	54	0.00000
536	54 PHE	H	-48.18040	-18.87361	116.41568	W.T	54	0.00000
537	54 PHE	CA	-46.18727	-19.64350	116.09999	W.I		0.00000
538	54 PHE	CB			116.63413	A1	54	0.00000
539	54 PHE	CC	-46.90625	-20.99411	118.13255		54	
540	54 PHE	CD1	-48.11656	-21.50316	118.65539	A1	54	0.00000
541	54 PHE	CD2	-45.89246	-20.53119	119.00621	Al.	54	0.00000
542	54 PHE	CEl	-48.31310	-21.55376	120.05266	A1	54	0.00000
543	54 PHE	CE2	-46.08993	-20.57973	120.40382	Al	54	0.00000
544	54 PHE	CZ	-47.30008	-21.09192	120.92307	Al	54	0.00000
545	54 PHE	С	-45.57270	-19.90650	114.73758	Al	54	0.00000
546	54 PHE	0	-44.36030	-19.93109	114.55246	ΥŢ	54	0.00000
5,47	55 GLU	N	-46.46681	-20.08976	113.76558	ΥŢ	55	0.00000
5,48	55 GLU	H.	-47.45338	-20.04226	113.95516	Al	55	0.00000
549	55 GLU	CA	-45.97322	-20.41545	112.42886	Al	55	0.00000
550	55 GLU	CB	-47.14512	-20.81915	111.54876	A1	55	0.00000
551	55 GLU	CG	-47.92567	-21.96126	112.19274	Al	55	0.00000
552	55 GLU	.CD	-49.04456	-22.36202		λl	55	0.00000
553	55 GLU	OE 1	-49.15553	-23.53977		Al	55	0.00000
554	55 GLU	OE2	-49.80524	-21.50984	110.81474	ΥŢ	55	0.00000
555	55 GLU	c:	-45.19354	-19.30188	111.76318	Al	55	0.00000
55.6	55 GLU	Ο.	-44.12678	-19.50387	111.19343	X1	55	0.00000
557	56 ALA	И	-45.73650	-18.08673	111.89532	Al	56	0.00000
558	56 ALA	H	-46.62336	-17.96765	112.34890	AI	56	0.00000
559	56 ALA	CA	-45.00414	-16.92733	111.38640	Al	56	0.00000
560	56 ALA	ÇВ	-45.80074	-15.63871	111.59969	Al	56	0.00000
561	56 ALA	C	-43.63772	-16.77849	112.02791	Al	56	0.00000
562	56 ALA	0	-42.62065	-16.60634	111.36878	Al	56	0.00000
563	57 GLN	N	-43.63088	-16.91454	113.35929	A1	57 57	0.00000
564	57 GLN	H	-44.48826	-17.03983	113.86832	ΥŢ	57	0.00000
565	57 GLN	СУ	-42.35063	-16.88200	114.06847	A1	57	0.00000
566	57 GLN	CB	-42.61987	-17.03049	115.56930	A1	57	0.00000
567	57 GLN	CG	-41.40659	-16.75355	116.45895	A.1	57	0.00000
566	57 GLN	CD	-41.80070	- i é . 94589	117.90800	A.1	57	0.00000
569	57 GLN	OEl	-42.29879	-16.06206	118.58910	Al	57	0.00000
570	57 GLN	NE2	-41.55425	-18.16011	116.38366	ā.l	57	0.00000
571	57 GLN	HE21	-41.14709	-18.87176	117.81428	Al	57	0.00000
572	57 GLN	HE22	-41.78115	-15.36728	119.33313	A.1	57	0.00000

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573	57 GLN	С	1.3621	317.94055	113.59025	λ1	57	0.0000
574	57 GLN	_	-40.1897	1 -17.68051	113.34773	Al	57	0.00000
373	58 GLY		-41.88825	5 -19.15742	113.41319	נע	္ 8 ၈	0.00000
576	58 GLY		-42,8477	1 -19.33514	113.6484/	VI VI	58 58	0.00000
577	58 GLY		-41.05373	9 -20.23393 4 -19.92781	112.8/3/4	r A	58	0.00000
578	58 GLY 58 GLY		-40,45534	-20.01021	111.30337	λ	58	0.00000
579 580	59 ALA	0 N	-41 34654	-19.52227	110.60018	A1	59	0.00000
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582	59 ALA		-40.88666	-19.14537	109.26181	Αl	59	0.00000
583	59 ALA	СВ	-42.07124	-1B.74086	; 108.30101	λl	59	0.00000
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588 589	60 LEU 60 LEU	CA CB	-39.17020 -39.87974	-14.72108	110.88234	A1	60	0.00000
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591	60 LEU	CDl	-41.72077	-13.05637	110.83175	Al	60	0.00000
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594	60 LEU	0	-36.81366	-15.71151	110.64266	<i>y</i> 2	61	0.00000
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607 608	62 ASN 62 ASN	HD21	-35.71613	-20.90910	106.04169	Al	62	0.00000
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790	80 THR	CG2		-15.72265 -14.32258	93.03840 Al	80	0.00000
	80 THR	C		-14.97644	92.99811 A1	80	0.00000
791 79 2	80 THR 81 PRO	0		-13.08096	93.61451 A1	81	0.00000
793	81 FRO	СD И		-12.25452	94.50529 A1	81	0.00000
794	81 PRO	CD CA		-12.43774	92.32255 A1	81	0.00000
795	81 PRO	CB		-11.39225	92.28996 A1	81	0.00000
796	81 PRO	CG		-10.94901	93.74456 21	81	0.00000
797	81 PRO	Č		-11.80328	92.25040 A1	81	0.00000
798	81 PRO	ρ.		-10.58936	92.16776 Al	81	Ü.00000
799	82 ILE	Ň .	-14.55883		92.27154 A1	82	0.00000
800	82 ILE	н		-13.66392	92.31920 A1	82	0.00000
801	82 ILE	CA		-12.18130	92.13921 Al	82	0.00000
802	82 ILE	CB	-12.60010	-11.87598	93.53959 Al	82	0.00000
803	82 ILE	CG2	-12.41140	-13.13773	94.38674 A1	82	0.00000
804	82 ILE	CG1	-11.31152		93.44331 Al	82	0.00000
805	82 ILE	CD		-10.62709	94.80896 A1	82	0.00000
806	82 ILE	С	-12.35649		91.37376 Al	82	0.00000
B 07	82 ILE	OCT1		-12.81747	90.75062 A1	82 82	0.00000 0.00000
808	82 ILE		-12.72556	-14.3/446	91.38671 Al 94.99084 Bl	1	0.00000
809	B3 GLY	N	-17.53322	-0.31236	94.35235 B1	ì	0.00000
810	83 GLY	HTl	-17.21994	-1.12219	94.96444 B1	1	0.00000
811 812	83 GLY	HT2	-16.86357 -17.61098	0.01920	95.97150 B1	î	0.00000
813	83 GLY	HT3 CA	•	-0.91116	94.55151 B1	ī	0.00000
814	83 GLY	C	-18.52573	-2.38203	94.66351 B1	1	0.00000
815	83 GTA	õ		-2.70920	94.84086 B1	1	0.00000
816	84 ASP	N	-19.57260	-3.20239	94.59303 Bl	2	0.00000
817	84 ASP	H .	-20.49658	-2.89510	94.35600 B1	2	0.00000
818	84 ASP	CA	-19.43900	-4.63200	94.86181 Bl	2	0.00000
819	84 ASP	СВ	-19.44643	-4.83356	96.38475 Bl	2	0.0000
820	84 ASP	CG	-18.89301	-6.19619	96.73815 Bl	2	0.00000
821	04 ASP	OD1	-17.69451	-6.29764	96.98109 Bl	2	0.00000
822	84 ASP	OD2	-19.66566	-7.14958	96.75715 B1	2	0.00000
923	84 ASP	С	-20.62766	-5.31072	94.19848 B1	2	0.00000
824	84 ASP	0	-21.46903	-4.61697	93.63479 B1	2	0.00000
925	85 THR	ĸ	-20.67796	-6.64606	94.24891 B1	3	0.00000
826	85 THR	н	-20.04362	-7.15926	94.84032 Bl	3 3	0.00000 0.00000
E27	65 THR	CH	-21.75257	-7.39367	93.59154 Ei	3	0.00000
328	95 THR	C5	-21.58903	-7.31950	92.05122 E1	2	0.0000

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829	85 TH	R OG1	32.7626	5 -7.82480	91.39964 B3	. 3	0.00000
830	85 TH		-22.7343			_	0.00000
83%	85 TH		-20.3296	6 -8.02078	91.53071 B	. 3	0.00000
832	85 TH		-21.74290	-8.83243	94.09035 B1		0.00000
833	85 TH		-20.76454			3	0.00000
834	86 AR		-22.86491	L -9.52779	93.88799 Bl		0.00000
835	B6 AR		-23.62404				0.00000
836	B6 AR			-10.87552			0.00000
837	86 ARC			-10.79083			0.00000
838	86 ARC			-11.65365			0.00000
839 840	86 ARC 86 ARC			-13.10294		4	0.00000
841	86 ARG			-14.03911 -14.68836		4	0.00000
842	86 ARG	-		-14.05016	· · · · · · · · · · · · · · · · · · ·	4	0.00000
843	86 ARG			-14.85617		4	0.00000
. 844	86 ARG		-19.73815		99.26299 Bl	4	0.00000
845	86 ARG	HH12	-20.05194	-15.47498	97.69021 Bl	4	0.00000
846	36 ARG			-13.26636	99.42563 B1	4	0.00000
847	86 ARG				100.26897 B1	4	0.00000
848 .	86 ARG				· 99.36831 B1	4	0.00000
849	86 ARG			-11.82065	93.57905 B1	4	0.00000
850 851	86 ARG			-11.52883		4	0.00000
852	87 PRO 87 PRO			-12.98532	93.27325 B1	- 5 5	0.00000
853	87 PRO			-13.34562 -14.02757	93.54376 B1 92.53087 B1	5	0.00000
854	87 PRO			-15.13066	92.39999 B1	5	0.00000
855	87 PRO			-14.43376	92.52460 B1	5	0.00000
856	87 PRO		-25.10390		93.21975 B1	5	0.00000
857	87 PRO	0	-25.32441		94.41838 B1	5	0.00000
858	88 ARG	N	-25.94344	-15.17061	· 92.39123 B1	6	0.00000
859	88 ARG	н	-25.68611	-15.35588	91.44433 B1	6	0.00000
8 60	88 ARG	CA	-27.23219	-15.62675	92.90235 Bl	6	0.00000
861	BP ARG	CB	-28.29078		91.81653 B1	6	0.00000
862	88 ARG	CG	-29.70863		92.35844 B1	6	0.00000
B 63	88 ARG	CD	-30.79150		91.49236 B1	6 6	0.00000
8 6 4 8 6 5	88 ARG	ne He	-31.92741 -31.76512		92.34607 B1 93.33788 B1	6	0.00000
866	88 ARG		-33.08881		91.86193 B1	6	0.00000
867	88 ARG		-34.06565		92.71322 B1	6	0.00000
868	88 ARG		-34.95441		92.40144 B1	6	0.00000
869	88 ARG	HH12	-33.91853	-13.92533	93.69838 B1	6	0.00000
870	88 ARG		-33.26623		90.54974 B1	6	0.00000
871	88 ARG		-34.12737		90.16725 Bl	6	0.00000
872 873	88 ARG		-32.52080		89.92214 B1 93.36366 B1	6 6	0.00000 0.00000
874	BB ARG BB ARG		-27.23157 -26.89591		92.63574 B1	6	0.00000
875	89 PHE		-27.62757 ·		94.62546 B1	7	0.00000
876	89 PHE		-27.92346		95.18173 B1	7	0.00000
877	89 PHE		-27.64368		95.17433 B1	7	0.00000
878	89 PHE		-26.56379 -	-18.74887	96.25427 B1	7	0.0000
879	89 PHZ	CG -	-25.20774 -	-18.55805	95.61829 B1	7	0.0000
880	89 PHE		-24.74915 -	-19.48139	94.64858 B1	7	0.00000
881	89 PHE		-24.42840 -		95.96102 B1	7	0.00000
882	89 PHE		-23.51045 -		94.00961 B1	7	0.00000
883 884	249 PKS		-23.18767 <i>-</i>		95.32384 B1	7	0.00000 0.00000
884 885	89 PHE 89 PHE		-22.73800 -		94.34976 B1 95.72084 B1	7	0.00000
986	89 PHE		-28.99632 - -29.85551 -	-10.3/14/ .18 12/29	95.94786 B1	, ל	0.00000
887	90 LED	й -	-29.85352 - -29.15755 -	20.28822	95.87791 B1	8	0.00000
886	90 LEU	H -	26.38720 -	20.92114	95.76376 B1	8	0.00000
889	90 LEU	CA ~	-30.48975 -	20.82226	96.14113 B1	3	0.00000
890	90 LEU	CB -	31.01265 -	21.46546	94.64981 B1	8	0.00000
891	90 LEU	CG -	32.46079 -	21.16892	94.43614 B1	ε	C.00000
292	90 LEU	CD1 -	32.81137 -	21.96230	93.19113 21	3	0.00000

FIG. 15

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893	90 LE	CD2	3.4749	8 -21.4069	9 95.55222	B1	8	0.00000
894	90 LEU			7 -21.9028			8	0.00000
895	90 LEU	_		6 -22.8920	-		В	0.00000
896	91 TRE			3 -21.7108			9.	0.00000
897	91 TRP			7 -20.8522			9	0.00000
898	91 TRP			7 -22.8604			9	0.00000
899	91 TRP		-31.7615	9 -22.5078	9 100.54418		9	0.00000
900	91 TRP		-30.4605	0 -22.1149	0 101.18157	Bl	9	0.00000
901	91 TRP			3 -22,7928		Bl	9	0.00000
902	91 TRP		-28.3007	4 -21.9654	6 101.98847	Bl	9	0.00000
903	91 TRP	CE3		B -24.0032		Bl	9	0.00000
904	91 TRP			0 -20.9642		Bl	9	0.00000
905	91 TRP			0 -20.8704	7 102.44030		9	0.00000
906	91 TRP	HE1		6 -20.1561		B1	9	0.00000
907	91 TRP	CZ2		7 -22.40520		Bl	9	0.00000
908	91 TRP	CZ3	-27.4405	4 -24.4040	100.81577	B1	9	0.00000
909	91 TRP	CH2		4 -23.61859	101.58828		9	0.00000
910	91 TRP	С	-32.96765	5 -23.39534	98.70806		9	0.00000
911	91 TRP	0		5 -22.6602			9	0.00000
912	92 GLN	N	33.02847	7 -24.71550	98.69197		10	0.00000
913	92 GLN	H		2 -25.28365			10	0.00000
914	92 GLN	CA		5 -25.34751			10	0.00000
915	92 GTN	CB		5,-,26.04119			10	0.00000
916	92 GLN	CG		-25.05724			10	0.00000
917	92 GLN	CD	-34.12668				10	0.00000
918	92 GLN	OE1		-25.88838			10	0.00000
919	92 GLN	NE2		-26.32196			10	0.00000
920	92 GLN	HE21					10	0.00000
921	92.GLN	HE22		<u>-26.83019</u>			10	0.00000
922	92 GLN	С		-26.35610	99.77327		10	0.00000
923	92 ĠTM	٥.	-33.40619	-27.07419		B1	10	0.00000
924	93 LEU	и.		-26.39205			11	0.00000
925	93 LEU	H	-36.24432	-25.74963	100.29272	D.I	11	0.00000
926	93 LEU	Cγ	-35.59559	-27.43909	101.48633	ם ז	11	0.00000
927	93 LEU	CB	-36.41178	-26.81504	102.88181	D1	11	0.00000
928	93 LEU 93 LEU	CG.	-34.835//	727.70766	103.99848	B1	11	0.00000
929 930	93 LEU	CD1 CD2	-35.82812			Bl	11	0.00000
931	93 LEU	CD2	-36 03506		101.32695		11	0.00000
932	93 LEU	Ö	-37.91692	-27 47571	100.96621	B1	11	0.00000
933	94 LYS	N.	-36 95045	-29 41719	101.56549	Bl	12	0.00000
934	94 LYS	н		-29.89066		B1	12	0.00000
935	94 LYS	CA		-30.14121		Bl	12	0.00000
936	94 LYS	·CB	-38.26049	-30.88530	100.13839	Bl	12	0.00000
937	94 LYS	CG	-39.43761	-31.85442	100.07544	Bl	12	0.00000
938	94 LYS	CD	-39.91969	-32.44832	98.75113	Bl	12	0.00000
939	94 LYS	CE	-41.00198	-33.40442	99.23924		12	0.00000
940	94 LYS	NZ		-33.90963	98.26846		12	0.0000
941	94 LYS	#21	-42.69798	-34.41995	98.86351		12	0.0000
942	94 LY8	HZ2		-33.14263	97.78947	B1	12	0.00000
943	94 LYS	HZ3	-41.57162	-34.57937	97.59540		12	0.00000
944	94 LYS	C	-38.34910	-31.08699	102.67642	Bl	12	0.00000
945	94 LYS	0	-37.54720	-31.99770	102.85849	B1	12	0.00000
946	95 PHE	N .	-39.40676	-30.85929	103.46609	B1	13	0.00000
947	95 PHE	H	-39.99321	-30.05201	103.34519	Bl	13	0.00000
948	95 PHE	CA	-39.73794	-31.84982	104.49053	R1	13	0.00000 0.00000
949	95 PHE	CB	-39.13251	-31.49228	105.86481	ET.	13	0.00000
950	95 PHE	CG	-39.62104	-30.20819	106.49960	D.1	13 13	0.00000
951	95 PHE	CD1	-39.C4917	-28.96886	106.12996	ים דם	13	0.00000
952	95 PHE	CD2	-40.60668	-30.26479	107.51173	D 1	13	0.00000
953	95 PHE	CE1	-39.45587	-27.78326	106.77985	# J	13	0.00000
954	95 PKE	CE2	-41.01477	-29.UBU19	108.16313	וע	13	0.00000
955 956	95 PHE	CZ	-40.4335/ -41.53005	-27.03433	104.57423	E)	13	0.00000
956	95 PHE	С	-41.22005	32.20/33				-

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957	95 PHE	•	~ .08591	-31.31332	104.41691	Bl	1'3	-05-00000
958	96 GLU); O	47812	-33.46734	104.75822	Bl	14	0.0000
9::9	50 CLU	н	-40.74093	-34.09071	105.01948		14	0.00000
960	96 GLU	CA	-42.80002	-33.99059	104.40078		1 4	0.00000
961	96 GLU	CB	-42.75085	-34.25375	102.88321		14	0.00000 0.00000
962	96 GLU	CG	-43.92070	~34.87965	102.11791		14 14	0.00000
963	96 GLU	CD	-43.55505	-34.92496	100.64018		14	0.00000
964	96 GW	OEl	-44.08203	-34.15253 -35.68559	100.23154		14	0.00000
965	96 GLU	OE2	-43.13129		105.17925		14	0.00000
966 967	96 GLÚ	C O	-42 27510	-36,11395	105.37879	Bl	14	0.00000
968	97 CYS	n	-44 39621	-35.34431	105.62501	Bl	15	0.0000
969	97 CYS	н	-45 05281	-34,60566	105.44206	BI	15	0.00000
970	97 CYS	CA	-44.78990	-36.60747	106.25700	BI	15	0.00000
971	97 CYS	CB	-45.09404	-36.46870	107.76554	Bl	15	0.00000 0.00000
972	97 CYS	SG	-46.49873		108.42288	BI	15 15	0.00000
973	97 CYS	С	-45.89248		105.55706	D1	15	0.00000
974	97 CYS	0	-46.75253	-36.80546	104.88787	Bl	16	0.00000
975	98 HIS	N		-38.69951 -39.11332	305.71704 306.26371·	Bl	16	0.00000
976	98-HIS	H	-45.06854		105.05785	B1	16	0.00000
977	98 HIS	CA	-46.75592 -46.00867	,	104.19437	81	16	0.00000
978 979	98 HIS 98 HIS	CB CG	-45.38253		102.97364	Bl	16	0.0000
980	98 HIS	NDI	-45.88967	-40.13539	101.74183	B 1	16	0.00000
981	98 HIS	HD1	-46.72307	-40.60043	101.51181	B1	16	0.00000
982	98 HIS	CD2	-44.21286	-39.22166	102.89430	B1	16	0.00000 0.00000
983	98 HIS	NE2	-44.02278		101.58753	BI	16 16	0.00000
984	98 HIS	CE1	-45.05268	. • • • • • •	100.87633		16	0.00000
985	98 HIS	C	-47.56750			B1	16	0.00000
986	98 HIS	0	-47.04279		105.81218	91	17	0.00000
987	99 PHE	N	-48.87987 -49.25243		105.00666	Bl	17	0.00000
. 988	99 PHE 99 PHÉ	H CA	-49.78029			Bl	17	0.00000
989 990	99 PHE	CB	-50.48946		107.59234	81	17	0.00000
991	99 PHE	CG	-49.63868	-39.72312	108.81497	Bl	17	0.00000
992	99 PHÉ	CD1	-48.50685	-38.87580	108.76543	B1	17	0.00000 0.00000
993	99 PHE	CD2	-49.92563	-40.47457	109.97439	Bl	17 17	0.00000
994	99 PHE	CE1	-47.64103		111.08354	B1	17	0.00000
995	99 PHE	CE2	-49.06252	-40.39779 -39.56658		Bl	17	0.00000
996	99 PHE	CZ	-47.92137		106.19800	Bl	17	0.0000
997	99 PHE	Ë	-50.79242 -51.48007		105.22101	Bl	17	0.00000
998 999	99 PHE	и 0	-50 86837	-43.12653	106.85844	Bl	18	0.00000
1000	100 PHE	.H	-SO 32768	-43.27017	107.68944	Bl	18	0.00000
1001	100 PHE	CA	-51 R4718	-44,10783	106.41132	81	18	0.00000 0.00000
1002	100 PHE	CB	-51.44468	-45.52210	106.84262	B1	18 18	0.00000
1003	100 PHE	CG	-51.08740	-46.34931	105.63001	8.1 P.T	18	0.00000
1004	100 PHE	CD1	-49.81159	-46.95065 -46.50158	103.33300	B1	18	0.0000
1005	100 PHE	CD2	-52.01269	-47.69723	104,30322	Bl	18	0.0000
1006	100 PHE	CE1	-49.45012	-47.24829	103.42719	Bl	18	0.00000
1007 1008	100 PHE	CE2 CZ	_50 37408	-47.83B59	103.35608	BI	18	0.00000
1008	100 PKE	C	-53 21727	-43.84401	106.96975	BI	18	0.00000
1010	100 PHE	0	-53 38235	-43.49692	108.13319	BI	18	0.00000 0.00000
1011	101 ASK	N	-54 19611	-44.08275	106.08672	27	19	0.00000
1012	101 ASN	н	-53 92777	-44.28429	105.14366	BI	19 19	0.00000
1013	101 ASN	CA	-55.63451	-43.97453	100.3/2/3	P.1	19	0.00000
1014	101 ASN	CB	-56.35400	-45.03359 -44.83624	105.52933	21	19	0.00000
1015	101 ASN	CG	-5/.66040	-44.83624	104.77523	21	19	0.00000
1016	101 ASK	0D1 ND2	_50 51 777	-45.52184	206.41359	81	19	0.00000
1017	101 ASN	ND21	_ ED DARAS	-46.21753	10/.02623	= 1	: 9	0.00000
1018 1015	101 ASN	HD22	EQ E0666	-05 49785	106.46862	= 7	19	0.00000
1019	101 ASX	С	-55.06277	-44.09262	107.83398	91	1 9	0.00000
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1021	10	l asn	0	16 73737	-43.2399	7 108.39463	Bl	19	0.00000
1022		2 GLY	-		-45.1894			20	0.00000
2025		2 GLY	• •		-45.0218			20	0.00000
1024		2 GLY				3 109.88157		20	0.00000
1025		2 GLY				2 110.81953		20	0.00000
1026		2 GLY			-45.1786			20	0.00000
1027			-	-54.84744	-43.2971	6 110.54648	Bl	21	0.00000
1028		3 THR		-55.35512	-42.9073	1 109.77435	Bl	21	0.00000
1029	10	3 THR	CA	-53.96549	-42.39542	2 111.30363	Bl	21	0.00000
1030		3 THR	CB	-54.75487	-41.8126	3 112.51011	B1	21	0.00000
1031	10	3 THŔ	0 G1	-54.22405	-40.5335	5 112.87483	B1	21	0.00000
1032	10:	3 THR	HGl	-54.81974	-40.0991	9 113.49221	Вl	21	0.00000
1033	10:	3 THR	CG2	-54.86925	-42.7243	3 113.73959	Bl	21	0.00000
1034	103	3 THR	С	-52.56983	-42.92386	5 111.66956	Вl	21	0.00000
1035	10:	THR	0	-51.96086	-42.59288	112.68707		21	0.00000
1036	104	GLU	N	-52.05837	-43.78433	110.78327	Bl	22	0.00000
1037		GLU	16	-52.50836	-43.94021	109.90068		22	0.00000
1038		GLU	CA	-50.80234	-44.42458	111.16198	B1	22	0.00000
1039		GLU	CB	-50.88647	-45.95191	111.05395	B1	22 22	0.00000 0.00000
1040				·51.1336B	-46.62525	112.41371	10 J	22	0.00000
1041		GLU	CD	-50.07267	-46.21410	113.43075	D.1	22	0.00000
1042		GLU	0E1			114.57130		22	0.00000
1043		GLU	OE2	-4889638 40. 55503	-40.13140	113.08359 110.45954	ום	22	0.00000
1044		GLÜ	C			109.34894		22	0.00000
1045	104	-	0	749.58628	44 04570	109.34074	BI	23	0.00000
1046	105		N	-48.46470	-44.059/8	111.19345	בם	23	0.00000
1047		ARG	H	-48.46820	-44.6/993	111.97754	נפ	23	0.00000
1048		ARG	CŸ	-47.23862	-43.305/8	110.79677	DI DI	23	0.00000
1049		ARĢ	CB	-45.43619				23	0.00000
1050		ARG	CG	-45.1116/	-42.37595	111.65081		23	0.00000
1051		ARG	CD	-43.89570	-43.22016 -42 48704	111.89499	BI	23	0.00000
1052 1053		ARG	NE HE	-42.03040	-42.40709	112.83941	Bl	23	0.00000
1055		ARG	CZ	-42.00777	-41 77685	110.93892	Bl	23	0.00000
1055		ARG	NH1	-41.02352	-41 00938	111.25165		23	0.00000
1056		ARG	HH11					23	0.00000
1057		ARG	нн12	-40 73208	-40.92747	112.21979		23	0.00000
1058		ARG	NH2	-42 47661	-41.82718	109.68286	Bl	23	0.00000
1059	-	ARG	HH21	-42.19810	-41.15340	109.00310	Bl	23	0.00000
1060		ARG	RH22		-42.57763	109.38941	Bl	23	0.00000
1061	105		Ċ	-46.47373	-44.13883	109.74376	Bl	23	0.00000
1062	105	ARG	0	-46.02105	-45.25363	109.97034	Bl	23	0.00000
1063	106		N	-46.32883	-43.51517	108.56827	Bl	24	0.00000
1064	106	VAL	H	-46.72693	-42.61352	108.37187	B1	24	0.00000
1065	106	VAL	CA	-45.53349	-44.24499	107.58352	B1	24	0.00000
1066		VAL	CB.	-46.27081	-44.39073	106.24298	B1	24	0.00000
1067		VAL	CG1	-45.79579	-45.65575	105.52366	B1	24	0.00000
1068		VIL	CG2	-47.77990	-44.41166	106.42922	Bl	24	0.00000 0.00000
1069		AVL	C	-44.14065	-43.66075	107.41554	RT.	24	0.00000
1070		VAL	0	-43.30354	-43.80891	108.29696	B1	24 25	0.00000
1071	107		И	-43.87314	-42.98069	106.29416	D.T	25	0.00000
1072	107		H	-44.58452	-42.69040	105.656/1	D.I	25	0.00000
1073	107		CA	-42.49561	-42.53115	106.12329	D1 D1	25	0.00000
1074		ARG	CB	-41.95685	-43.01317	104.76032	D 3	25	0.00000
1075	107		CG	-41.96328 -42.33357	-52.U2619 -12 66620	102.30304	R1	25	0.00000
1076	107		CD	-42.33357 -43.76838	-42.00029 -23 63515	102.27063	E1	25	0.00000
1077	107		NE 	-44.37946	-42.33313	102.43147	B1	25	0.00000
1078			HE CZ	-44.26510		102.20945	B1	25	0.0000
1079 1080	107		NH1	-45.56598	-44 37747	102.35905	5.1 B.1	25	000000
1081	107		PH11	-46.08629	-45.19707	102.34562	51	25	0.00000
1081	107		HH12	-45.19539	-43.55730	102.54612	E 3	25	0.00000
1093	107		11:12	-43.47983	-45.23695	102.05793	5 1	25	0.00000
1084	10.		HH21	-43.86541 ·	-46.15096	102.05452	21	25	0.00000

_لمعدر.	LTN2 . CPD		<u> ಇಗಿಬ್ಗಳಿಕರಿ</u>	25 14:58:4	8 1993		10	
1085	107 ARG	нн22	2.49404	-45.11997	101.94496	Bl	25	0.00000
1086	107 ARG	С	2.37787	-41.03546	106.32504	Bl	25	0.00000
1007	157 ARG	0	-43.36896	-40.31981	106.43645	Bl	25	0.00000
1088	108 LEU	И	-41.12633	-40.58602	106.39627		26	0.00000
1089	108 LEU	H	- 40.32617	-41.17985	106.29320	B1	26 26	0.00000 0.00000
1090	108 LEU		-40.89942	-39.15744	106.53751		26	0.00000
1091	108 LEU	CB	-40.80087	-38.80954 -37.89689	108.01240		26	0.00000
1092	108 LEU	CG	-41.90347 -42.75820	-38.44729			26	0.00000
1093 1094	108 LEU	CD1	-41 50070	-36.44650			26	0.00000
1094	108 LEU	C	-39 62061	-38,77273	105.85764	Вl	26	0.00000
1096	108 LEU	õ	-38 76368	-39,61093	105.60154	Bl	26	0.00000
1097	109 LEU	N	-39.52150	-37.47828	105.56773	Bl	27	0.00000
1098	109 LEU	н	-40 28459	-36.84167	105.72467	B1	27	0.00000
1099	109 .TER	CA	-38.30683	-37.02679	104,90334	B1	27 27	0.00000 0.00000
1100	109 LEU	CB	-38.47418	-37.24039	103.39009	BJ RT	27	0.00000
1101	109 LEU	CG	-39.74490	-36.59695 -35.25631	102.82839	ופ	27	0.00000
1102	109 LEU	CD1	-39.42210	-35.25631 -37.56523	101 91138	Bl	27	0.00000
1103	109 LEU 109-LEU	CD2	-40.49190	-35.58061	-105.21268	B1	27	0.00000
1104 1105	109 LEU	0	-34 AR411	-34.77322	105.47943	Bl	27	0.00000
1105	110 GLU	N	-36,70092	-35.29768	105.12126	Bl	28	0.00000
1107	110 GLU	11	-36.01498	36.01786	105.00992	Bl	28	0.00000
1108	110 GLU	CV	-36.20315	-33.92985	105.08184	Вl	28	0.00000
1109	110 GLU	CB	-35.80977	-33.47627	106.49578	Bl	28	0.00000
1110	110 GLÚ	CĢ	-35.32216	-32.02210	106.60993	B1	28	0.00000
1111	110 GLU	CD		-31.87178	106.30413	B1	28	0.00000 0.00000
1112	110 GLU	OE1	-33.37957	-30.74951	106.10521	B1	28 28	0.00000
1113	110 GLU	OE2	-33.09671	-32.85201	106.36491	D1	28	0.00000
1114	110 GLU	C	-35.01026	-33.94076	104.15500	B1	28	0.00000
1115	110 GFA	0	-34.25577	-34.90807	104.11363	Bl	29	0.00000
1116	111 ARG	N	-34.89439	-32.86843 -32.16766	103.37222	Bl	29	0.00000
1117	111 ARG	Н Са	-33.01363	-32.74903	102.46249	B1	29	0.0000
1118 1119	111 ARG	CB	-33.99004	-33.60407	101.20352	B1	29	0.00000
1120	111 ARG	CG	-35.45106	-33.70037	100.75305	Bl	29	0.00000
1121	111 ARG	CD	-35.67880	-34.83036	99.75174	Bl	29	0.00000
1122	111 ARG	NE	-37.09124	-35.21147	99.72614	Bl	29	0.00000 0.00000
1123	111 ARG	HE	-37.71860	-34.61357	100.22757	B1	29 29	0.00000
1124	111 ARG	CZ	-37.50098	-36.32384	99.09789 99.16119		29	0.00000
1125	111 ARG	NH1	-38.78116	-36.689/I	98.68276	Bl	29	0.00000
1126	111 ARG	HH11	-39.12269 -39.43712	-36 16699	99.70940	B1	29	0.00000
1127 1128	111 ARG 111 ARG	NH2	-36.63391	-37 06893	98.41518	Bl	29	0.00000
1129	111 ARG	ин21	-36.91501	-37.90348	97.94325	Bl	29	0.00000
1130	111 ARG	нн22	-35.67503	-36.78688	.98.36264	Bl	29	0.00000
1131	111 ARG	С	-33 49135	-31.31194	102.08170	Bl	29	0.00000
1132	111 ARG	0	-34.39395	-30.54569	101.76599	31	29	0.00000
1133	112 CYS	N	-32.20568	-30.97057	102.12414	B1	30	0.00000
1134	112 CYS	н.	-31.50733	-31.64380	102.37455	B1	30	0.00000 0.00000
1135	112 CYŞ	CA	-31.80468	-29.62360	101.73826	B1	30 30	0.00000
1136	112 CYS	CB	-31.12874 -30.70297	-28.92303	102.51550	Bl	30	0.00000
1137	112 CYS	SG	-30.70297	-29 64998	100.54380	Bl	30	0.00000
1138 1139	112 CYS 112 CYS	С 0	-29.97769	-30.48052	100.40406	Bl	30	0.00000
1140	112 LIS	Ŋ	-31.15975	-28.70696	99.65078	BI	31	0.00000
1141	113 ILE	н	-31.84193	-28.00116	99.86111	81	31	0.00000
1142	113 ILE	CA.	-30.55306	-28.70228	98.32464	Bl	31	0.00000
1143	113 ILE	СЭ	-31.56021	-29.33526	97.30706	Bl	31	0.00000
1144	113 ILE	CG2	-33.00339	-29.37434	97.82323	B.7	31	0.00000
1145	113 ILE	CG1	-31.52431	-28.74184	95.89635	ום	31 31	0.00000
1146	113 ILE	CD	-32.44047	-29.50025	94.93272 97.9£536	B) PT	31	0.00000
1147	113 ILE	C	-30.08576	-21.29000	98.21317	B1	31	0.00000
1148	113 ILE	C	-30.75333	-29.29900	30.2.20-1			

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1149	114 2	YR N	28.8701	0 -27.2556	4 97.41665 E	1 32	0.0000
1150				0 -28.1015			0.00000
1151	114 2	YR CA	-28.2203	6 -25.9826			0.00000
1152	114 2	YR CB	-26.8008	8 -26.0193			0.00000
1153				7 -24.7210			0.00000
1154				0 -23.4582			0.00000
1155				6 -22.2808			0.00000
1156				4 -24.8021			0.00000
1157				1 -23.6265			0.00000
1158	114 7			0 -22.3694			0.00000
1159		YR OH		7 -21.2191° 6 -21.43015			0.00000
1160 1161	114 7			-21.43013 -25.8321			0.00000
1162	114 T			-26.70965			0.00000
1163	115 A			-24.6999			0.00000
1164	115 A			-24.02391			0.00000
1165	115 A			-24.41545	93.68566 B	1 33	0.00000
1166	115 A	SN CB	-27.27049	-23.84385			0.00000
1167	115 A			-22.43118			0.00000
1168					94.85370 B		0.00000
1169	115 አ			-21.48359		_	0.00000
1170	115 A		-28.04883				0.00000
1171	115 A		-2729503	. •			0.00000
1172	115 አ			-25.59300			0.00000
1173	115 A			-25.91402 -26.23583			0.00000
1174	116 G			-25.92614			0.00000
1175 1176	116 G			-27.38897			0.00000
1177	116 G			-26.98635			0.00000
1178	116 G		-31.33401	-27.96419	89.80983 B		0.00000
1179	116 G			-27.31473	88.43588 B	34	0.00000
1180	116 G			-27.33002			0.00000
1181	116 G		-30.20428	-26.71450			0.00000
1182	116 G			-26.71396			0.00000
1183	116 G1			-26.24641	87.20205 B		0.00000
1184	116 GI			-28.73663			0.00000
1185	116 GI			-29.77893	92.10767 B 93.24838 B		0.00000
1186	117 GI 117 GI			-28.72803 -27.88576	93.59525 B		0.00000
1187 1188	117 GI			-30.02538	93.58636 B		0.00000
1189	117 GI			-30.02385	93.23572 B		0.00000
1190	117 GI		-26.41125	-29.71932	91.75724 B		0.0000
1191	117 GI		-24.93459	-29.87167	91.43636 B		0.00000
1192	117 GI	•	-24.62388	-30.40776	90.37359 B	35	0.00000
1193	117 GL		-24.09945	-29.45965	92.24135 B	35	0.00000
1194	117 GL				95.05360 B		0.00000
1195	117 GL		-28.39032		95.93593 B: 95.31151 B:		0.00000 0.00000
1196	118 GL	•	-28.45418		94.57056 B		0.00000
1197 1198	118 GF		-28.37884 -28.64640		96.70504 B	36	0.00000
1199	118 GL		-28.86529		96.80220 B		0.00000
1200	118 GL		-30.04821		96.04454 B		0.00000
1201	118 GL		-30.18685		96.39969 B	36	0.0000
1202	118 G		-31.31366		96.62004 B	36	0.00000
1203	118 GL	Ú OE2	-29.17775	-36.44658	96.46236 B		0.00000
1204	118 GL	U C.	-27.45968		97.59954 B		0.00000
1205	118 GE	•	-26.30375		97.22004 B		0.00000
1206	119 SE		-27.77719		98.81671 B		0.00000
1207	119 SE		-28.73032	-31.22468	99.10568 53		0.00000
1208	119 SE		-26.67523	-31.09310	99.74333 B		0.00000
1209	119 SE	R CB	-26.79679	-27.04425 -20 22240	100.25438 B3 100.96854 B3	37	0.00000
1210	119 SE		-23.02500	-28 33730	101.30354 5	37	0.00000
1211	119 SE		-25.75540	-32.10869	100.87899 B1	37	0.00005
			20.04507				

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1213	119 SER	٥	25.72568	-32.90554	101.03211	в1 3	
1213	120 VAL	Й	-27.72434	-32.06808	101.67561	B1 3	
1215	120 VAL	H	-28.50026	-31.45938	101.48344	B1 3	
1216	120 VAL	CA	-27.82678	-32.99802	102.80099	B1 3	
1217	120 VAL	CB	-27.44790	-32.32244	104.13888	B1 3	·
1218	120 VAL	CG1	-25.93252	-32.18078 -30.97900	104.29388		_
1219	120 VAL	CG2	-28.15631	-33.53625	107.33300		
1220	120 VAL 120 VAL	С 0	-29.23/1/	-32.88656	102.49711	B1 3	
1221	120 VAL	N	-29 34164	-34.75356	103.41944	B1 3	9 0.00000
1223	121 ARG	н	~28.52525	-35.27675	103.67737	B 1 3	
1224	121 ARG	CA	-30.65941	-35.37282	103.49447	B1 3	
1225	121 ARG	CB	-30,83879	-36.26556	102.25451	B1 3	
1226	121 ARG	CG	-32.24891	-36.83907	102.13977	B1 3	
1227	121 ARG	CD	-32.36448	-38.11228	101.31436	B1 3 B1 3	
1228	121 ARG	NE	-33.46636	-38.89614	101.86585	B1 3	=
1229	121 ARG	HE	-33.58314	-38.82444 -39.71705	102.88363		
1230	121 ARG	CZ	-34.21779	-40.40193	101.73316	B1 3	
1231	121 ARG	NH1	-35.77971	-41.03741	.101.23780	B1 3	
1232 1233	121 ARG	HH12	-35.33365	-40.28453	102.71808	в1 3	9 0.00000
1233	121 ARG	NH2	-33.99786	-39.84536	99.82444	B1 3	
1235	121 ARG	HH21	-34,53079	40.47140	99.25675		
1236	121 ARG	HH22	-33.27054	-39.30635	99.39731	B1 3	
1237	121 ARG	С	-30.80743	-36.21013	104.75602	B1 3	
1238	121 ARG	0	-29.86515	-36.82872	105.23483	B1 3	
1239	122 PHE	N	-32.04075	-36.24095 -35.63810	103.20404	B1 4	
1240	122 PHE	H	-32.75061 -32.40668	-37.26064	106.24769	B1 4	
1241	122 PHE 122 PHE	CA CB	-32.40000	-36.87165	106.85296	B1 4	0.00000
1242	122 PHE	CG	-33.64992	-36.46831	108.30418	B1 4	
1244	122 PHE	CDI	-32.56100	-35.69712	108.77664	B1 4	
1245	122 PHE	CD2	-34.66985	-36.87929	109.19094	B1 4	
1246	122 PHE	CEl	-32.49275	-35.34303	110.14035	B1 4	
1247	122 PHÉ	CE2	-34.60187	-36.52254	110.55431	B1 4	
1248	122 PHC	CZ	-33.51285	-35.75823	111.02379	B1 4	
1249	122 PHE	С	-32.57844	-38.62424 -38.79208	103,60697		•
1250	122 PHE	0	-33.34100	-39.60796	106.15389	B1 4	
1251 1252	123 ASP 123 ASP	N H	-31.23184	-39.47015	106.92548	B1 4	0.00000
1253	123 ASP	CA	-32.08552	-40.93522	105.58825	B1 4	
1254	123 ASP	СВ	-30.85171	-41.81726	105.76445	BI 4	
1255	123 ASP	CG	-29.93161	-41.56041	104.59405	B1 4	
1256	123 ASP	001	-28.81173	-41.11310	104.81757		
1257	123 ASP	OD2	-30.34905	-41.80191 -41.63618	105.43838	B1 4	
1258	123 ASP	C	-33.32362	-41.03010	107.02319	B1 4	
1259 1260	123 ASP 124 SER	о И	-33 62443	-42.74250	105.41429	B1 4	
1261	124 SER	Н	-32.95329	-43.10971	104.76900	B1 4	
1262	124 SER	CA.	-34.94425	-43.35498	105.58880	B1 4.	
1263	124 SER	СВ	-35.18779	-44.37028	104.46149	B1 4	
1264	124 SER	OG	-36.57945	-44.70776	104.37671	B1 4:	
1265	124 SER	HG	-36.91089	-44.87590	105.27413	B1 4	
1266	124 SER	С	-35.21640	-44.01057 -44.42223	107.22372		
1267	124 SER	0	-30.33338	-44.10325	107.74330	B1 4	
1268	125 ASP	N H	-33.2622A	-43.73425	107.50494	B1 4	0.00000
1269 1270	125 ASP 125 ASP	CA	-34 30492	-44.60689	109.10471	B1 4	
1271	125 ASP	CB	-32.96210	-45.24580	109.50620	B1 4	
1272	125 ASP	CG	-31 82155	-44.23245	109.55779	P1 4	
1273	125 ASP	001	-31 R0194	-43.19553	108.89075	21 4	
1274	125 ASP	OD2	-30 87249	-44.45676	110.29614	B1 4	-
1275	125 ASP	С	-34.65374	-43.52244	110.10428		
1276	125 ASP	0	-35.10540	-43.78441	111.22672		

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1277	126 VAL	и	51615	5 -42.26751	109.64790	B1 4	4 0.00000
1278	126 VAL	н	2093 دن-	-42.12308	108.70456	B1 4	4 0.00000
1279	126 VAL	СУ	-34.59346	6 -41.07914	. 110.50496	B1 4	4 0.00000
1280	126 VAL	CB	-36.04814	4 -40.82462	110.97592	B1 4	4 0.00000
1281	126 VAL	CG1	-36.22164	-39.41758	111.54094	B1 4	4 0.00000
1282	126 VAL	CG2	-37.04494	-41.002B1	109.82481		4 0.00000
1283	126 VAL	С	-33.57625	5 -41.13739	111.64961		4 0.00000
1284	126 VAL	. 0	-33.78819	-40.75401	112.79533		4 0.00000
1285	127 GLY	N	-32.41547	-41.66443	111.25324		5 0.00000
1286	127 GLY	H		-42.01122			5 0.00000
1287	127 GLY	CA		-41.88201		B1 4	
1288	127 GLY	С		-41.56889	111.62542	B1 4	
1289	127 GLY	0	-29.03055			B1 4	
1290	128 GLU	N	-29.81143	-41.65704		B1 4	
1291	128 GLU	H		-42.13586 -41.14297		B1 4	
1292	128 GLU 128 GLU	CA CB		-42.29481			
1293 1294	128 GLU	CG		-42.25781			6 0.00000
1295	128 GLU	CD		-43.28422			6 0.00000
1296	128 -GLU	OE1-	24.49646	-42.88073			6 0.00000
1297	128 GLU	OE2		-44.47773		Bl 4	
1290	128 GLU	C	-28.71376	-39.98012	108.75031	B1 4	
1299	128 GLU	ō		39.57375		B1 4	6 0.00000
1300	129 TYR	N	-27.54735	-39.42619	108.38931	B1 4	0.00000
1301	129 TYR	н	-26.68498	-39.83420	108.68820	B1 4	
1302	129 TYŔ	CA.	-27.50019	-38.29454	107.46434	B1 4	
1303	129 TYR	CB	-26.63842	-37.15326	108.01560		
1304	129 TYR	CG	-27.30857	-36.34222	109.09443		
1305	129 TYR	CDļ		-36.21958			
1306	129 TYR	CE1		-35.42131			
1307	129 TYR	CD2	-28.52827	-35.67832	108.83210		
1308	129 TYR	CE2		-34.88063	109.83571		
1309	129 TYR	CZ	-28.47327			B1 4	
1310	129 TYR	OH		-33.98459 -33.40059		B1 4	
1311	129 TYR	нн	-29.70394 -26.82531		106.15941	B1 4	
1312 1313	129 TYR 129 TYR	C		-39.04407		B1 4	
1313	129 TYR 130 ARG	0	-27.55686		105.08581		B 0.00000
1315	130 ARG	н		-38.10451	105.16735	B1 4	B 0.00000
1316	130 ARG	CA		-38.41254	103.80227	B1 4	
1317	130 ARG	CB		-38.84699	102.71647		
1318	130 ARG	CG	-27.21143	-39.00526	101.34112		
1319	130 ARG	CD		-39.41974	100.29808	B1 4	
1320	130 ARG	'NE		-39.38333	98.95629	41	
1321	130 ARG	HE		-38.85034	98.82833		
1322	130 ARG	CZ	-28.29934	-39.96348	97.93202		
1323	130 ARG	ИНЛ		-39.81313	96.69917 ,95.90996		
1324	130 ARG	HHII	-28.26738	-40.23513 -39.25297	96.53955		
1325 1326	130 ARG			-40.68730	98.14392	B1 4	
1327	130 ARG 130 ARG	NH2		-41.10492	97.39118		B 0.00000
1328	130 ARG			-40.81930	99.07957		0.0000
1329	130 ARG	C		-37.06053	103.45986		0.00000
1330	130 ARG	ŏ		-36.11772	103.07688		
1331	131 ALA	ĸ	-24.95816	-36.99899	103.58668	B1 4	
1332	131 ALA	н	-24.45069	-37.77908	103.95192	B1 4	
1333	131 ALA	CA	-24.28607	-35.84894	102.98902	B1 4	
1334	131 ALA	СВ	-23.06137	-35.44271	103.80801	51 49	
1335	131 ALA	С	-23.85084	-36.20633	101.58658		
1336	131 ALA	0		-37.20532	101.36065		
1337	132 VAL	N		-35.36774	100.63154	B1 50	
1338	132 VAL	li O		~34.54695	100.85567	B1 50	
1339	132 VAL	CA	-24.05930	-35.79619	98.36560		
1340	132 VAL	CB	-25.125/8	-35.11200	50.50500	<u> </u>	

./DRl_	MIN2 . CP	5	Tou Feb	25 14:58:4	8 1993	22	
1341	132 VA	L CG1	5.02505	5 -35.45687	96.87606		
1342	132 VA		-26.51048	3 -35.48853	98.89777		
1243	132 VX			-35.53393	98.73419		
1344	132 VA			3 -36.19257	97.84114	B1 50	
1345	133 TH	R N	22.01646	5 -34.53567	99.35299		
1346	133 TH	R H	-22.44036	-34.04272	100.11190	B1 51	
1347	133 TH	R CA		-34.13851	98.91128		
1348	133 TH	R CB	-20.84979	-33.09962	97.76600		0.00000
1349	133 TH	•	-19.58519	-32.56706	97.34397		0.00000
1350	133 TH		-19.69018	-32.10136	96.50551 98.12788		0.00000
1351	133 TH		-21.81730	-31.96947			
1352	133 TH		-19.91735	-33.59750 -33.30952	100.10840	B1 51	0.00000
1353	133 TH		-20.48697	-33.30952 -33.44216	99.91793	B1 52	0.00000
1354	134 GL		-18.60113	-33.65725	99.02398		0.00000
1355	134 GL		-18.20332	-32.88738	100.97647		0.00000
1356	134 GLU		-17,75230		100.49482		0.00000
1357	134 GLU		-15.30332		100.03844		0.00000
1358 1359	134 GLU		-14 31377	-34.04410	99.60488		0.00000
1360	134-614	I - OEI-		-34.96158			0.00000
1361	134 GLU		-13.96021	-33.03519	98.99487	B1 52	0.00000
1362	134 GLU		-18.18601	-31.51310	101.45728	B1 52	0.00000
1363	134 GLU		-17.97884	.,-31.11916	102.59352	B1 52	0.00000
1364	135 LEU		-18.87611	-30.80674	100.55812	B1 53	0.00000
1365	135 LEU		-18.92762	-31.13340	99.61398	B1 53	0.00000
1366	135 LEU	CA	-19.55727	-29.57029	100.94931	B1 53	0.00000
1367	135 LEC	Г СВ	-20.29914		99.70959	B1 53	0.00000
1368	135 LEU	CG	-20.12211	-27.57937	99.38843		0.00000
1369	135 LEU	CD1	-20.93390	-26.68772	100.32666	B1 53	0.00000
1370	135 LEU	CD2	-18.63204	-27.23870	.99.36436		0.00000
1371	135 LEU	C		-29.72915	102.11412		0.00000
1372	135 LEU		-20.60025	-28.93514			0.00000
1373	136 GLY		-21.29082	-30.82221	102.02891	B1 54	0.00000
1374	136 GLY		-21.14062	-31.49656	101.30480	B1 54	0.00000
1375	136 GLY		-22.25373	-31.10000	104.26834		0.00000
1376	136 GLY		-21.66447	-31.85099 -31.82013	105 38053	B1 54	0.00000
1377	136 GLY 137 ARG		-24,17,60	-32.53644	104.01306	B1 55	0.00000
1378 1379	137 ARG		-20.34132	-32.57514			0.00000
1380	137 ARG		-19.94165	-33.32437			0.0000
1381	137 ARG			-34.02762		B1 55	0.00000
1382	137 ARG		-18.13008	-35.07893			0.00000
1383	137 ARG		-19.16855	-36.15719	105.85371	B1 55	0.00000
1384	137 ARG		-18.66781	-37.06640	106.87657	B1 55	0.00000
1385	137 ARG	HE	-17.93002	-36.71980	107.47204	B1 55	0.00000
1386	137 ARG		-19.21896	-38.26612	107.07086	B 1 55	0.00000
1387	137 ARG	NH1	-18.67934	-39.07643	107.97489	B1 55	0.00000
1388	137 ARG	HH11	-19.03969	-39.99255	108.14749	B1 55	0.00000
1389	137 ARG	HH12	-17.88763	-38.75541	108.49972	B1 55	0.00000
1390	137 ARG	NH2	-20.29085	-38.64369	106.37219	B1 55 B1 55	0.00000
1391	137 ARG		-20.72192	-39.53705	105.49834	B1 55	0.00000
1392	137 ARG	нн22	-20.68879	-38.01490 -32.65026	105.70437	B1 55	0.0000
1393	137 ARG	Ç,	-19.76330	-33.13638	107 44172	B1 55	0.0000
1394	137 ARG 138 PRO	0 N	-10 03005	-31.51331	106.56277		0.00000
1395	138 PRO	CD	-18.26903	-30.78028	105.55482		0.00000
1396 1397	138 PRO	CA	-18 92500	-30.86885	107.87839	B1 56	0.00000
1398	138 PRO	СВ	-18 00966	-29.66775	107.61418	51 56	0.00000
1398	138 PRO	CG	-17.22921	-30.01427	106.35389	B1 56	0.00000
1500	138 PRO	C	-20.24853	-30.39451	108.45452	<u>21 56</u>	0.00000
1401	138 PRO	õ	-20.38873	-30.20105	109.65243	31 56	0.00000
1402	139 ASP	N	-21,22822	-30.19487	107.56513	31 21	0.00000
1403	139 ASR	H	-21.12955	-30.45258	106.60295	21 57	0.00000
1404	139 ASF	CA	-22.54445	-29.76374	108.02635	B1 57	0.00000

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1405	139 ASP	СВ	3.29481	-29.2279	4 106.80729	B1	5 7·	~0~0 00 000
1406					107.23406		57	0.00000
1,07					5 107.53698		57	0.00000
1408			-25,59590	-29.0093	7 107.24891	Bl	57	0.00000
1409			23.29009	-30.9065	L 108.70646	Bl	57	0.00000
1410		_	-23.84428	-30.7994	109.79628	Bl	57	0.00000
1411	140 ALA		-23.18867	-32.06950	108.04975	Bl	58	0.00000
1412			-22.77745	-32.09325	5 107.13514	Bl	58	0.00000
1413			-23.64141	-33.29863	108.70194	Bl	58	0.00000
1414	140 ALA		-23.39932	-34.51285	107.80148	Bl	58	0.00000
1415	140 ALA		-22.96994	-33.52593	110.04660	B1	58	0.00000
1416	140 ALA		-23.61501	-33.62460	111.08086	Bl	58	0.00000
1417	141 GLU		-21.63141	-33.53710	110.01537	Bl	59	0.00000
1418	141 GLU	Н	-21.14260	-33.44422	109.14251	Bl	59	0.00000
1419	141 GLU	CA	-20.88131	-33.73045	111.26272	Bl	59	0.00000
1420	141 GLU	СВ	-19.38545	-33.75474	110.92637	Bl	59	0.00000
1421	141 GLU	CG	-19.08157	-34.92155	109.97183	Bl	59	0.00000
1422	141 GLU	CD	-17.65605	-34.91070	109.44677	Bl	59	0.00000
1423	141 GLU	OE1	-17.21662	-35.95752	108.96182	Bl	59	0.00000
1424	141 GLU	· OE2	·· -16.9965B	-33.87375	-109.50288	Bl	59	0.00000
1425	141 GLU	С	-21.20315	-32.72395	112.36696	Bl	59	0.00000
1426	141 GLU	0	-21.35204	-33.05368	113.53869	Bl	59	0.00000
1427	142 TYR	N	-21.39109	31.47026	.111.93865	в1	60	0.00000
1428	142 TYR	H	-21.16858	-31.22939	110.99159	Bl	60	0.00000
1429	142 TYR	CA	-21.91640	-30.42572	112.82625	Вl	60	0.00000
1430	142 TYR	CB	-22.17510	-29.18770	111.95478	Bl	60	0.00000
1431	142 TYR	CG	-22.15441	-27.86866	112.68902	Bl	60	0.00000
1432	142 TYR	CD1	-20.91930	-27.21121	112.08286	Вl	60	0.00000
1433	142 ŤYŘ	CE1	-20.89216	-25.94177	113.49633		60	0.00000
1434	142 TYR	CD2		-27.27306			60	0.00000
1435	142 TYR	CE2	-23.33600	-26.00211	113.72688		60	0.00000
1436	142 TYR	CZ	-22.10013	-25.34007	113.91274	-	60	0.00000
1437	142 TYR	ОН	-22.06472	-24.08718	114.49226		60	0.00000
1438	142 TYR	нн	-22.95958	-23.75980	114.62492	B1	60	0.00000
1439	142 TYR	Ç	-23.20365	-30.84932	113.52485	B1	60	0.00000
1440	142 TYR	0	-23.33185	-30.86335	114.74427	B1	60	0.00000
1441	143 TRP	N	-24.16819	-31.24530	112.69102	E1	61	0.00000
1442	143 TRP	H	-24.01203	-31.27353	111.69669	B1	61	0.00000
1443	143 TRR	CA	-25.46084	-31.65772	113.24427	B1	61	0.00000
2444	143 TRP	CB	-26.46502	-31.82534	112.10045	B.T	61 61	0.00000
1445	143 TRP	CG	-26.82927	-30.51319	111.43167	D.T	61	0.00000
1446	143 TRP	CD2	-27.59514	-30.35383	110.26190 110.01923	ום	61	0.00000
1447	143 TRP	CE2	-27.68725	-28.88192	109.37526	BI	61	0.00000
1448 1449	143 TRP 143 TRP	CE3	-26.23171	-20 2040B	111.84528	Bl	61	0.00000
1450	143 TRP	CD1 NE1	-20.43303 -25 00373	-28 24161	111.01939	Bl	61	0.00000
1451	143 TRP	HE1	-26 86335	-27 27405	111.09738	B1	61	0.00000
1452	143 TRP	CZZ	-28 41151	-28.41903	108.90296	Bl	61	0.00000
1453	143 TRP	CZ3	-28 94655	-30.73482	108.27096	B1	61	0.00000
1454	143 TRP	CH2	-29 03488	-29.34388	108.03833	Bl	61	0.0000
1455	143 TRP	C .	-25 40824	-32.93379	114.07770	Bl	61	0.00000
1456	143 TRP	0	-26 13451	-33.11650	115.04995	B1	61	0.00000
1457	144 ABN	N	-24 46546	-33,80055	113.69236	Bl	62	0.00000
1458	144 ASN	Ж М	-23 04027	-33.62305	112.85783	B1	62	0.00000
1459	144 ASN	CA	-24.16067	-34,99080	114.49069	Bl	62	0.00000
1460	144 ASN	CB	-23.20850	-35.93308	113.73882	B1	62	0.00000
1451	144 ASN	CG	-23.89541	-36.68861	112.61740	Bl	62	0.00000
1462	144 ASN	ODI	-23.68155	-36.47075	111.43371	Bl	62	0.00000
1463	144 ASN	ND2	-24.72776	-37.64097	113.02365	B1	62	0.00000
1464	144 ASN	HD21	-24.89338	-37.61926	113.9927€	Bl	62	G.C0000
1465	144 ASN	HD22	-25 19584	-39.20406	112.34455	B 1	62	0.00000
1466	144 ASN	Ç	-23 49875	-34.69497	115.82591	В1	62	0.00000
1 4 5 7	144 ASN	0	-23.43003	-35.54654	116.69934	91	62	o.00000
1468	145 SER	1:	-22.99604	-33.46640	115.97353	31	63	0.00000

.,		ن:	عدد دهد	. LD 14:55:	ذولات ٥٠	24	
1469			22.9783	0 -32.8042	8 115.21961 11	69 🛀	.00000:0
1470		R CA	22.3800	4 -33.1462	1 117.25865 B1	63	0.00000
3.47.3		_			4 117.01176 B1	63	0.0000
1472					0 118.24511 B1	63	0.00000
1473					5 118.91065 B1	63	0.00000
1474					4 118.06161 B1	63	.0.00000
1475					7 119.13941 B1	63	0.00000
1476					7 117.51252 B1	64	0.00000
1477	146 GLN				116.64774 B1	64	0.00000
1478	146 GLN				5 118.19522 B1	64	0.00000
1479	146 GLN				117.23892 B1	64	0.00000
1480	146 GLN	-			117.72882 B1	64	0.00000
1481	146 GLN				2 117.25234 B1	64 64	0.00000
1482 1483	146 GLN 146 GLN				118.00911 B1 115.94329 B1	64	0.00000
1484	146 GLN				115.33276 B1	64	0.00000
1485	146 GLN				115.56134 B1	64	0.00000
1486	146 GLN		-26.29488	3 -30.93694	118.57642 B1	64	0.00000
1487	146 GLN	ō			117.77294 B1	64	0.00000
1488	147 LYS				119.82371 B1	65	0.00000
1489	147 LYS	н			120.47663 B1	65	0.00000
1490	147 LYS	CA	-27.35463	32.46614	120.21964 B1	65	0.00000
1491	147 LYS	CB	-28.27258	32.00851	121.37895 B1	65	0.00000
1492	147 LYS	CG			121.84203 B1	65	0.00000
1493	147 LYS	CD			121.97244 B1	£ 5	0.0000
1494	147 LYS	CE			121.47444 B1	65	0.00000
1495	147 LYS	NZ			121.26907 Bl	€5	0.00000
1496	147 LYS	H21		-37.42223		65	0.00000
1497	147 LYS	H22			120.68922 B1	65	0.00000
1498	147 LYS	HZ3			122.15073 B1	65	0.00000
1499	147 LYS	C			119.09340 Bl	65	0.00000
1500	147 LYS	0		-32.90174	118.83883 B1	65	0.00000
1501	148 ASP	N		-34.00916	118.45822 B1 118.75300 B1	66 66	0.00000 0.00000
1502	148 ASP	H		-34.14495 -35.00236	117.48362 B1	66	0.00000
1503 1504	148 ASP 148 ASP	CA CB		-36.37833		66	0.00000
1505	148 ASR	CG		-36.57930	119.21585 B1	66	0.00000
1506	148 ASP	OD1	-25.69280	-36.06325		66	0.00000
1507	148 ASP	OD2		-37.20742	120.23524 B1	66	0.00000
1508	148 ASP	c .		-34.63043	116.79645 B1	66	0.00000
1509	148 ASP	0 -		-35.14808		66	0.0000
1510	149 LEU	14	-28.92399	-33.61840	115.93941 B1	67	0.00000
1511	149 LEU	H		-33.30949		67	0.00000
1512	149 LEŲ	· CA		-32.84836	115.45008 B1	67	0.00000
1513	149 LEU	CB	-29.45399	-31.74267	114.57360 B1	67	0.00000
1514	149 LEU	CG	-30.29432	-30.58223	114.02475 B1	67	0.00000
1515	149 LEU	CD1	-30.85820	-30.92475	112.65290 Bl	67	0.00000 0.00000
1516	149 LEU	CD2	-31.34761	-30.09615	115.02072 B1	67 67	0.00000
1517	149 LEU	C			114.80952 B1	67 67	0.00000
1518 1519	149 LEU	0	-32.36472	-33.3/839	114.83807 Bl 114.32363 Bl	68	0.00000
1520	150 LEU	н У	-20.75110	-34.00130	114.11621 B1	68	0.00000
1521	150 LEU	CA			114.12850 Bl	68	0.00000
1522	150 LEU	CB			114.15286 B1	68	0.00000
1523	150 LEU	CG			112.89569 B1	68	0.00000
1524	150 LEU		-28.64060	-38.26626	113.18796 B1	68	0.00000
	150 LEU	CD2	-30.71709	-38.01712	111.73915 B1	68	0.00000
1526	150 LEU	С	-32.74973	-36.17247	115.10785 Bl	65	0.00000
	150 LEU	0	-33.89001	-36.01901	114.70350 Bl	63	0.00000
1528	151 GLU	ĸ	-32.47441	-36.43576	116.39428 B1	6.5	0.00000
	151 GLU	Н	-31.52943	-36.57284	116.72119 B1	65	0.00000
1530	151 GLU	CA	-33.61295	-36.59512	117.30950 El	69	0.00000
	151 GLU	CB	-33,19489	-36.98331	118.72928 B1	69	0.00000
1532	121 CTA	CG	-32.69081	-38.41906	118.86324 51	6 2	0.00000

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253					64 118.70497 E		0.0000
135					29 117.58148 B		0.00000
153					62 119.72459 B		0.00000
153					68 117.43857 B		0.00000
153					13 117.58647 B		0.00000
1539 1539		•			00 117.35586 B3		0.00000
1540				6 -33.047			0.00000
1543					6 117.32018 BI		0.00000
1542					00 118.60094 B1		0.00000
1543	152 GL				4 118.49387 B		0.00000
1544			-32.1306	8 -29.2701	.9 118.36008 B1		0.00000
1545					2 118.54711 B1		0.00000
1546					1 118.65663 B1		0.00000
1547					8 118.46911 B1		0.00000
1548 1549					6 116.27182 B1	70 70	0.00000
1550		-	,		1 116.43607 B1 9 115.08552 B1	70 71	0.00000
1551	153 ARG				9 114.99299 B1	71	0.00000
1552					9 119.95947 B1	71	0.00000
1553	153 ARG				7 112.66564 B1	71	0.00000
1554	153 ARG				1 112.35054 B1	71	0.00000
1555	153 ARG				0.111.17839 B1	71	0.00000
1556	153 ARG				6 110.55166 B1	71	0.00000
1557	153 ARG				0 110.98860 B1	71	0.00000
1558	153 APG				2 109.31716 B1	71	0.00000
1559	153 ARG				6 108.54980 B1	71	0.00000
1560	153 ARG	•			2 107.55422 B1	71	0.00000
1561	153 ARG				108.90136 Bl	71	0.00000
1562 1563	153 ARG 153 ARG				1 108.83735 B1	71	0.00000
1564	153 ARG 153 ARG	HH21 -	32.50269	-32.29042	2 107.04355 B1 9 109.42307 B1	71 71	0.00000 0.00000
1565	153 ARG				114.06926 B1	71	0.00000
1566	153 ARG				113.96948 B1	71	0.00000
1567	154 ARG				114.34589 B1	72	0.00000
1568	154 ARG				114.34685 B1	72	0.00000
1569	154 ARG	CA -:	37.32441	-37.02834	114.59374 B1	72	0.00000
1570	154 ARG				115.12926 B1	72	0.00000
1571	154 ARG				115.22908 B1	72	0.00000
1572	154 ARG				115.69023 B1	72	0.00000
1573 1574	154 ARG 154 ARG				115.80184 B1	72 72	0.00000
1575	154 ARG 154 ARG				115.09038 B1 116.80504 B1	72	0.00000 0.00000
1576	154 ARG				116.88687 B1	72	0.00000
1577	154 ARG	-			117.61587 B1	72	0.00000
1578	154 ARG	HH12 -3	8.17495	-43.96809	116.20728 B1	72	0.00000
1579	154 ARG	NH2 -3	5.75363	-42.48325	117.71513 B1	72	0.00000
1580	154 ARG				118.48012 B1	72	0.00000
1581	154 ARG				117.63709 B1	72	0.0000
1582	154 ARG				115.51516 B1	72	0.00000
1583 1584	154 ARG				115.28503 B1	72	0.00000
1585	155 ARG 155 ARG				116.55472 B1 116.77047 B1	73 73	0.00000 0.00000
	155 ARG				116.77047 B1	73	0.00000
	155 ARG				117.38371 B1	73	0.00000
	155 ARG				119.44775 B1	73	0.00000
	155 ARG				120.46128 31	73	0.00000
1590	155 ARG	NE -41	0.71623 -	-32.83780	121.28237 B1	. 73	0.00000
	155 ARG				121.09517 81	73	0.00000
	155 ARG				122.22743 91	7.5	0.00000
	155 ARG	NH1 -43	1.48382 -	-31.43599	122.94081 B1	73	0.00000
	155 ARG				123.65592 El	73 73	0.00000
	155 ARG	#### ~ 43	2.41818 - 3.99916	-31./5018	122.77039 B1 122.45304 B1	7.7 7.7	0.00000 0.00000
- 3 7 0	155 ARG	NH2 -39	= . 22 / 15 -	31.49569	122.43304 5:		0.00000

. ,	_مسارك . فتق)	Thu Feb	عد 14:58:	c8 1993	27	
1661	162 ARG	: 11E	50.4969	4 -40.88455	5 114.11092	B1 &0	J.000TO
1,662	162 ARG		49.3426	7 -42.4050	7 114.80601	B1 80	
70.		NH1	-50.2302	0 -43.32564	114.46730	B1 80	
1664	162 ARG	нні	1 -50.1197	9 -44.3140	114.66150	B1 80	
1665			2 -51.0618	4 -43.06283 0 -42 76163	5 113.96028 3 115.38323	B1 80	
1666 1667	162 ARG 162 ARG	NH2	-48.1963 1 -48 0013	4 -43.72773	115.55650	B1 80	
1668	162 ARG	HH2	2 -47.5257	9 -42.06788	115.64161	B1 80	0.00000
1669	162 ARG		-49.75512	2 -35.92906	5 113.42487		
1670	162 ARG	•	-50.82093	3 -35.93732	112.81658		
1671	163 HIS		-49.58593	3 -35.31862	114.60088		0.00000 0.00000
1672	163 HIS	H	-48.68252	2 -35.28035	115.03933 115.23467	B1 81	
1673 1674	163 HIS 163 HIS	CA CB	~50.76433 ~50.42874	-34.73337	116.66134	B1 81	0.00000
1675	163 HIS	CG	-50.41026	-35.50649	117.55950	B1 81	0.00000
1676	163 HIS	ND1	-51.51644	-36.01842	118.12240	B1 81	0.00000
1677	163 HIS	HD1	-52.42830	-35.67431	118.03242	B1 81	0.00000
1678	·163 HIS	CD2	-49.31404	-36.28837	117.93420	B1 81 B1 81	0.00000
1679 1680	163 HIS	NE2	-49.78030) -37.28089 -37.33585	118.73130	B1 81	0.00000
1681	163 HIS	C	-51.37160	-33.59588	114.44466	B1 81	0.00000
1682	163 HIS	ō	-52.56947	-33.55956	114.18659	B1 81	0.00000
1683	164 ASN	N	-50.50246	-32.67616	114.01202	B1 82	0.00000
1684	164 ASN	н		-32.74462			0.00000
1685	164 ASN	CA		-31.58239			0.00000
1686	164 ASN	CB	-49.96583	-30.54361 -29.53869	112.89196		0.00000
1687 1688	164 ASN 164 ASN	CG.	-49.91907	-29.62233			0.00000
1689	164 ASN	ND2		-28.55321			0.00000
1690	164 ASN			-28.49843			0.00000
1691	164 ASN	HD22	-50.84210	27.84178	114.61251	B1 82	0.00000
1692	164 ASN	C	-51.70674	-32.02333	111.91584	B1 82	0.00000
1693	164 ASN	0	-52.73418	-31.48662	111.51085	B1 82	0.00000
2694	165 TYR	N	-51.12326	-33.05626	111.29191	B1 83 B1 83	0.00000
1695	165 TYR	н	-50.28201	-33.49370	111.62194 110.10401	B1 83	0.00000
1696 1697	165 TYR 165 TYR	CA CB	-51.81304 -50.92781		109.25048	B1 83	0.00000
1698	165 TYR	CG	-51.39689	-34.50284	107.80317	B1 83	0.00000
1699	165 TYR	CD1	-52.18946	-33.44669	107.29137	B1 83	0.00000
1700	165 TYR	CEl	-52.63312	-33.47413	105.96126	B1 83	0.00000
1701	165 TYR	CD2	-51.04289	-35.58963	106.96903	B1 83 B1 83	0.00000
1702 1703	165 TYR 165 TYR	CE2 CZ	-51.48846	-34.55457	105.62740	B1 83	0.00000
1703	165 TYR		-52.75931	-34.53152			0.00000
1705	165 TYR	нн	-52.13515	-34.97162	103.24416	B1 83	0.00000
1706	165 TYR	Ċ	-53.16114	-34.17050	110.39688	B1 83	0.00000
1707	165 TYR	0	-54.17243	-33.75040	109.85354	B1 83	0.00000 0.00000
1708	166 GLY	N	-53.16827	-35.13255	111.32677	B1 84 B1 84	0.00000
1709	166 GLY	H	~52.30920	-35.55284	111.74443	B1 84	0.00000
1710 1711	166 GLY	CA C	-54.44388	-34 78683	112.08191	B1 84	0.00000
1712	166 GLY	0	-56.70058	-34.86763	111.64884	B1 84	0.00000
1713	167 VAL	N	-55.16433	-33.81049	112.90998	B1 85	0.00000
1714	167 VAL	H .	-54.21931	-33.77699	113.25020	B1 85	0.00000
1715	167 VAL	CA	-56.14083	-32.77588	113.26808	B1 85 B1 85	0.00000
1716	167 VAL	CB	-55.54422	-31.87200	114.36486	B1 85	0.00000
1717 1718	167 VAL 167 VAL	CG1 CG2	-55 22649	-30.71030	115.60849		0.00000
1719	167 VAL	CG2	-56.62003	-31.96437	112.06515	B1 85	.0.0000
1720	167 VAL	õ	-57.80658	-31.70971	111.87365	E1 85	0.00000
1721	168 GLY	N	-55.65605	-31.61592	111.20320	61 86	0.00000
1722	168 GLY	H	-54.68827	-31.81924	111.38362	B1 86	0.00000
:723	158 GLY	CA	-56.04215	-30,96490	109.94972	B1 86 B1 86	0.00000
1724	168 GLY	С	-57.02338	-31.//629	209.42907	2. 00	

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1725	168 GLY	. 0	58.02649	-31.28044	108.61617	Bl	86	0.00000
1726					109.00665		87	0.00000
1717			-55.87921				87	0.00000
1728			-57.59179	-34.00467	108.30469	Bl.	87	0.00000
1729	169 GLU	CB	-56.95070	-35.39155	108.28846	B1	87	0.00000
1730	169 GLU	CG	-55.67851				87	0.00000
1731			-54.91259				87	0.00000
1732	169 GLU	OE1	-53.74095				87	0.00000
1733	169 GLU	OE2	-55.48540				87	0.00000
1734	169 GLU	С	-59.00151				87	0.00000
1735	169 GLU	0	-59.98966	-34.06690	108.14126	B1	87	0.00000
1736			-59.06996	-33.99305	110.19884	B1	88	0.00000
1737			-58.23138	-34.05998	110.74831	B1	8 B 8 B	0.00000
1738	170 SER		-60.38255	-33.85094	110.83391	D.I	88	0.00000
1739	170 SER		-60.18950 ·	-33.03501	112.33790	B1	88	0.00000
1740 1741	170 SER 170 SER	OG HG	-61.30877				88	0.00000
1742	170 SER	C	-61.16415				88	0.00000
1743	170 SER	õ	-62.31497	-32.69150	109.96191	B1	88	0.00000
1744	171 PHE						89	0.00000
1745	171 PHE	н	-59.54601 -				89	0.00000
1746	171 PHE	CA	-61.19539 -				89	0.00000
1747	171 PHE	CB	-60.30793 -	-28.99941	110.10880	Bl	89	0.00000
1748	171 PHE	CC	-59.94208 -	-28.68147	111.54294	Bl	89	0.00000
1749	171 PHE	CD1	-58.59291 -	-28.39703	111.85413	Bl	89	0.00000
1750	171 PHE	CD2	-60.93098 -				89	0.00000
1751	171 PHÉ	CEl	-58.23032 -		113.17656		89	0.00000
1752	171 PHE	CE2	-60.56845 -		113.87858		89	0.00000
1753	171 PHĖ	CZ	-59.21901 -	-28.01319	114.18527		89	0.00000
1754	171 PHE	С	-61.62802 -		108.49502	B1	89	0.00000
1755	171 PHE	Ο,	-62.68697 -		108.07691		89 90	0.00000
1756	172 THR	N,	-60.72520 -		107.69903 108.09636		90	0.00000
1757 1758	172 THR 172 THR	H CA	-59.91792 - -60.84308 -				90	0.00000
1759	172 THR	CB	-59.41710 -	30.70747	105.70165	B1	90	0.00000
1760	172 THR	OG1	-59.35292 -		104.36207		90	0.00000
1761	172 THR	HG1	-60.15430 -		103.89285	Bl	90	0.00000
1762	172 THR	CG2	-58.90262 -	32.13906	105.74825	Bl	90	0.00000
1763	172 THR	Ç	-61.71208 -	31.64877	105.47987	Bl	90	0.00000
1764	172 THR	0	∺61.76078 -		104.25169		90	0.00000
1765	173 VAL	N	-62.32957 -		106.19315		91	0.00000
1766	173 VAL	H	-62.28706 -		107.19363		91	0.00000
1767	173 VAL	CA	-62.87984 -		105.53718		91 91	0.00000 0.00000
1768	173 VAL	. CB	-63.87967 - -64.50006 -		106.47899		91	0.00000
1769 1770	173 VAL	CG1	-63.20452 -	35.//324	107.23034	B)	91	0.00000
1771	173 VAL 173 VAL	CG2 C	-63.51710 -	34.90246	107.75400	a1	91	0.00000
1772	173 VAL	0	-63.18750 -		103.18126		91	0.00000
1773	174 GLN	N	-64.41211 -		104.14711		92	0.0000
1774	174 GLN	н	-64.59818 -				92	0.00000
1775	174 GLN	CA	-65.14373 -	32.19104	102.94243	B 1	92	0.00000
1776	174 GLN	Ca	-65.76132 -	30.80951	103.22574	B 1	92	0.0000
1777	174 GLN	CG	-66.77986 -	30.26016	102.21318	Bl	92	0.00000
1778	174 GLN	CD	-66.09397 -		101.04012		92	0.00000
1779	174 GLN	OE1	-65.45924 -		101.15380		92	0.00000
1780	174 GLN	NE2	-66.26402 -	30.1B961	99.87453		92 92	0.00000 0.00000
1781	174 GLN	HE21	-66.63226 -	31.11966	99.82928		92 92	0.00000
1782	174 GLN		-65.97833 -	29.14839 22.16624	99.02730 101.63553		92 92	0.00000
1783	174 GLN	C	-64.36067 - -64.88582 -		101.63333	Bl	92	0.00000
1784	174 GLN 175 ARG	0	-64.88582 -		101.73327		93	0.00000
1785 1786	175 ARG	H	-62.69754 -3		102.62225	Bì	93	0.0000
1787	175 ARG	C.A.	-62.31707 -	31.78835	100.50018	B1	93	0.00000
1788	175 ARG	CB	-61.85817 -	30.36631	100.16958	31	93	0.0000
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1789	175 ARG	CG	-61 2439	5 -30.26914	98.77310 B1	93	0.00000
1790	175 ARG	CD		7 -28.04209		93	0.00000
1791	175 ARG	NE		7 -28.87216		93	0.00000
1792	175 ARG	3H		6 -29.76910		93	0.00000
1793	175 ARG	CZ		5 -27.74924		93	0.00000
1794	175 ARG	NH1		3 -27.83579		93	0.00000
1795	175 ARG			4 -27.02728		93	0.00000
1796	175 ARG	HHI	2 -58.2248	0 -28.73459	95.46034 Bl	93	0.00000
1797	175 ARG	NH2	-59.9242	7 -26.55682	97.09593 B1	93	0.00000
1798	175 ARG	HH2:	-59.57021	8 -25.70796	96.70808 B1	93	0.00000
1799	175 ARG			1 -26.50975		93	0.00000
1800	175 ARG	C		6 -32.79150		93	0.00000
1891	175 ARG	0	-61.02327	7 -33.52738	99.49809 Bl	93	0.00000
1802	176 ARG	N	-60.36493	3 -32.85103	101.53216 B1	94	0.00000
1803	176 ARG	н	~60.53761	-32.31340	102.35642 B1	94	0.00000
1804	176 ARG	CA	-59.25084	-33.80802	101.41499 Bl	94	0.00000
1805	176 ARG	CB	-58.18750	-33.62483	102.49829 Bl	94	0.00000
1806	176 ARG	CG	-57.32871	-32.37173		94	0.00000
1807	176 ARG	CĐ	-56.27515	-32.28237	103.45854 Bl	94	0.00000
1808	176 ARG	NE	-56.30062	-30.94018	104.04511 B1	94	0.0000
1609	176 ARG	HΞ		-30.25818		94	0.0000
1810	176 ARG	CZ		-30.63659		94	0.00000
1811	176 ARG	NHI	1	29.38340		94	0.00000
1812	176 ARG	нніі				94	0.00000
1813	176 ARG	HH12				94	0.00000
1814	176 ARG	NH2			105.89297 B1	94	0.00000
1815	176 ARG	HH21				94	0.00000
1816	176 ARG	нн22		-32.51329	105.54660 B1	94	0.00000
1817	176 ARG	С	-59.64686			94	0.00000 0.00000
1818	176 ARG	0			101.06767 B1	94 95	0.00000
1819	177 VAL	N	-60.88626	-35.53390	101.85272 B1	95	0.00000
1820	177 VAL	H			102.18430 Bl 101.70299 Bl	95	0.00000
1621 1822	177 VAL	CA CB .		-36.91608 -37.47193		95	0.00000
1823	177 VAL	CB .		-39.00335		95	0.00000
1824	177 VAL	CG2		-37.00753		95	0.00000
1825	177 VAL	C		-37.02458	100.63553 B1	95	0.00000
1826	177 VAL	ō		-37.67295	100.79118 B1	95	0.00000
1827	178 HIS	N		-36.33028	99.52296 B1	96	0.00000
1828	178 HIS	н		-35.77698	99.39184 B1	96	0.00000
1829	178 HIS	CA		-36.32813	98.41659 B1	96	0.00000
1830	178 HIS	CB		-35.10064	98.57314 B1	96	0.00000
1831	178 HIS	CG		-35.21481	97.74669 B1	96	0.00000
1832		NDI		-35.62002	96.46944 Bl	96	0.00000
1833	178 HIS	HD1	- '	-35.91855	95.96816 Bl	96	0.00000
1834	178 HIS	CD2		-34.92566	98.14872 Bl	96	0.00000
1835	178 HIS	NE2		-35.16505	97.08780 Bl	96	0.00000
1836	178 HIS	CE1		-35.59364	96.05191 Bl	96	0.00000
1837	178 HIS	С	-62.34621	-36.29235	97.09131 Bl	96	0.00000
1038	178 HIS		-61.22615	-35.78756	97.07091 Bl	96	0.00000
1839	178 HIS	OCT2	-62.87363	-36.77415	96.08799 Bl	96	0.00000

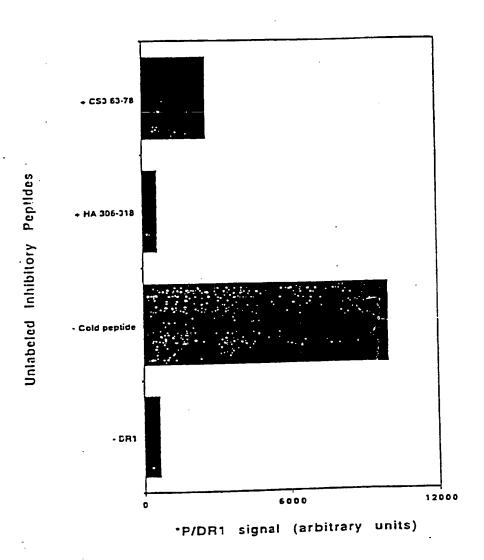
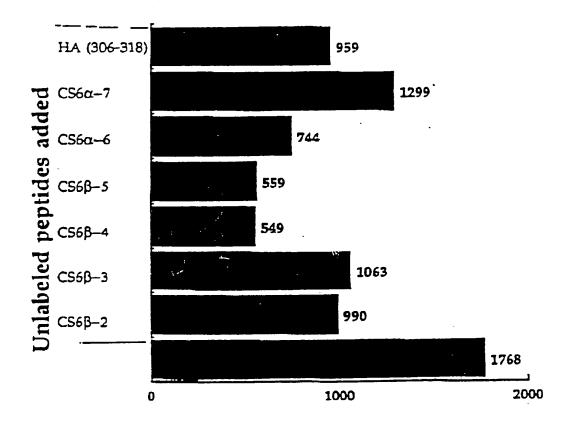


FIG. 31

Inhibition of 125 I HA (306-318)/DRI. by unlabeled CSG of and B peptides

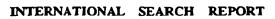


*HA/DR1 compact dimer signal (densitometric units)

INTERNATIONAL SEARCH REPORT

International application No.
PCT/US94/05697

US CL :	US CL: 424/185.1, 186.1, 190.1, 242.1; 530/327, 326, 333, 334, 388.75 According to International Patent Classification (IPC) or to both national classification and IPC							
	ocumentation searched (classification system followed	by classification symbols)						
1								
0.5. : 4	424/185.1, 186.1, 190.1, 242.1; 530/327, 326, 333,	334, 366.73						
Documentat	Documentation searched other than minimum documentation to the extent that such documents are included in the fields searched							
Electronic data base consulted during the international search (name of data base and, where practicable, search terms used)								
C. DOC	C. DOCUMENTS CONSIDERED TO BE RELEVANT							
Category*	Citation of document, with indication, where ap	propriate, of the relevant passages	Relevant to claim No.					
×	The Journal of Immunology, Volusius and 15 April 1993, Nauss et al. Peptides in a Structural Homology MHC ", page 41A, Abstract 221,	1, 3-20						
x	Nature, Volume 358, issued 27 A	uguet 1992 Chicz et al	12					
<u>^</u>	"Predominant Naturally Processed	•						
_Y	DR1 are derived from MHC-re		1, 3-7					
'	Heterogenous in Size", pages 764-							
	2, and Table 3.	, co, coo paga (ac, 1221)						
	2, 3							
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X Furth	er documents are listed in the continuation of Box C	See patent family annex.						
• Sp	ecial categories of cited documents:	"T" later document published after the inte date and not in conflict with the applic						
"A" document defining the general state of the art which is not considered to be of particular relevance		principle or theory underlying the inv						
	lier document published on or after the international filing date	"X" document of particular relevance; the considered novel or cannot be considered.						
"I." document which may throw doubts on priority claim(s) or which is cited to establish the publication date of another citation or other		when the document is taken alone	·					
special reason (as specified)		"Y" document of particular relevance; the considered to involve an inventive	step when the document is					
	cument referring to an oral disclosure, use, exhibition or other	combined with one or more other suc being obvious to a person skilled in t						
the	cument published prior to the international filing date but later than priority date claimed	*&* document member of the same patent family						
Date of the	actual completion of the international search	Date of mailing of the international sec	arch report					
O1 SEPTE	EMBER 1994	1 3 SEP 1994.						
Name and mailing address f the ISA/US Authorized officer			Do					
Box PCT	ner of Patents and Trademarks	H. Sidberry	za fa					
Washington Facsimile N	a, D.C. 20231	Telephone No. (703) 308-0196	/					
T seconime 14	o. (703) 305-3230	1.22,022,000						



International application No. PCT/US94/05697

Category*	Citation of document, with indication, where appropriate, of the relevant passages	Relevant to claim No
Y	The Journal of Immunology, Volume 150, No. 2, issued 15 January 1993, Boehncke et al., "The Importance of Dominant Negative Effects of Amino Acid Side Chain Substitution in Peptide-MHC Molecule Interactions and T Cell Recognition", pages 331-341, see Abstract, on page 331.	8-11
X.	The EMBO Journal, Volume 9, No. 6, issued 1990, Jardetzky et al., "Peptide binding to HLA-DR1: a Peptide with most residues substituted to alanine retains MHC binding", pages 1797-1803, see page 1798, page 1800, figure 4, and page 1801, figure 7.	512
Ÿ	Nature, Volume 332, issued 28 April 1988, Brown et al., "A hypothetical model of the foreign antigen binding site of Class II histocompatibility molecules", pages 845-850, see pages 845-849.	1, 3, 4
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Form PCT/ISA/210 (continuation of second sheet)(July 1992)*

INTERNATIONAL SEARCH REPORT

International application No.
PCT/US94/05697

Box 1 Observations where certain claims were found unsearchable (Continuation of item 1 f first sheet)					
This international report has not been established in respect of certain claims under Article 17(2)(a) for the following reasons:					
Claims Nos.: 2 because they relate to subject matter not required to be searched by this Authority, namely:					
Claim 2 is directed to a computerized model which encompasses scientific theory and computer programs to the extent that the International Searching Authority is not equipped to search prior art concerning such programs. Accordingly claim 2 is withdrawn from search under PCT Rule 39 and PCT Article 17(2)(a)(i).					
Claims Nos.: because they relate to parts of the international application that do not comply with the prescribed requirements to such an extent that no meaningful international search can be carried out, specifically:					
3. Claims Nos.: because they are dependent claims and are not drafted in accordance with the second and third sentences of Rule 6.4(a).					
Box II Observations where unity of invention is lacking (Continuation of item 2 of first sheet)					
This International Searching Authority found multiple inventions in this international application, as follows:					
1. As all required additional search fees were timely paid by the applicant, this international search report covers all searchable claims.					
2. As all searchable claims could be searched without effort justifying an additional fee, this Authority did not invite payment of any additional fee.					
3. As only some of the required additional search fees were timely paid by the applicant, this international search report covers only those claims for which fees were paid, specifically claims Nos.:					
4. No required additional search fees were timely paid by the applicant. Consequently, this international search report is restricted to the invention first mentioned in the claims; it is covered by claims Nos.:					
Remark n Protest The additional search fees were accompanied by the applicant's protest. N protest accompanied the payment of additional search fees.					

Form PCT/ISA/210 (continuation of first sheet(1))(July 1992)*